



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 10:51 AM GMT

PDB ID : 3MK3  
Title : Crystal structure of Lumazine synthase from Salmonella typhimurium LT2  
Authors : Kumar, P.; Singh, M.; Karthikeyan, S.  
Deposited on : 2010-04-14  
Resolution : 3.57 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.  
We welcome your comments at [validation@mail.wwpdb.org](mailto:validation@mail.wwpdb.org)  
A user guide is available at  
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>  
with specific help available everywhere you see the ⓘ symbol.

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The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467  
Mogul : 1.7 (RC4), CSD as536be (2015)  
Xtriage (Phenix) : 1.9-1692  
EDS : rb-20026688  
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)  
Refmac : 5.8.0135  
CCP4 : 6.5.0  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : trunk26865

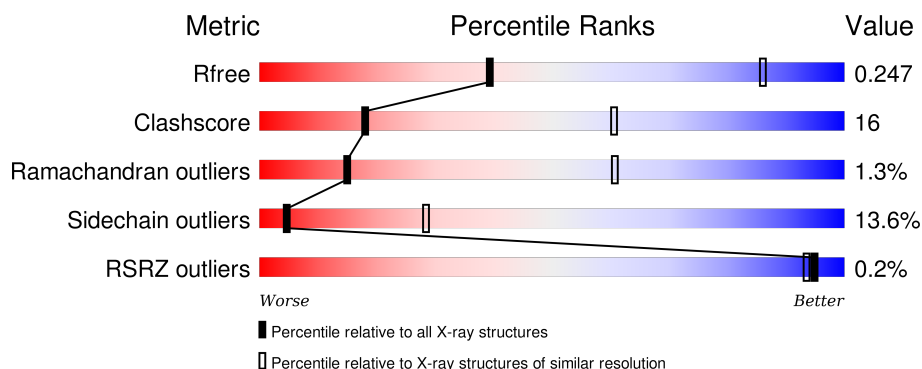
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*






The reported resolution of this entry is 3.57 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.








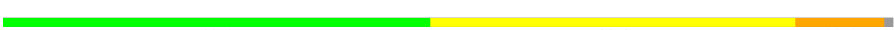




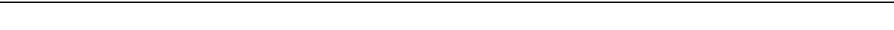

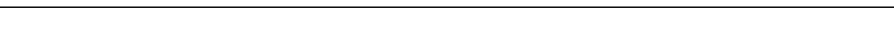
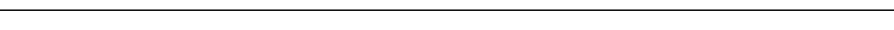











Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
$R_{free}$	91344	1240 (3.72-3.40)
Clashscore	102246	1057 (3.70-3.42)
Ramachandran outliers	100387	1017 (3.70-3.42)
Sidechain outliers	100360	1017 (3.70-3.42)
RSRZ outliers	91569	1247 (3.72-3.40)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions  $\leq 5\%$ . The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	1	156	 60% 30% 8% .
1	2	156	 60% 30% 8% .
1	3	156	 60% 30% 9% .
1	4	156	 58% 31% 10% .
1	5	156	 57% 33% 9% .











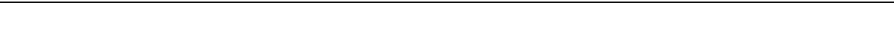

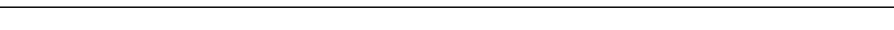
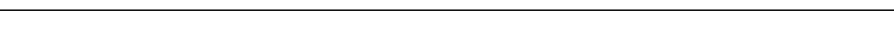











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Mol	Chain	Length	Quality of chain
1	6	156	 59% 32% 8% .
1	7	156	 58% 33% 8% .
1	8	156	 58% 31% 9% .
1	9	156	 60% 30% 9% .
1	A	156	 56% 34% 8% .
1	B	156	 48% 41% 10% .
1	C	156	 54% 35% 9% .
1	D	156	 54% 35% 9% .
1	E	156	 53% 37% 9% .
1	F	156	 49% 42% 8% .
1	G	156	 51% 40% 8% .
1	H	156	 52% 37% 10% .
1	I	156	 58% 34% 7% .
1	J	156	 54% 35% 10% .
1	K	156	 54% 36% 9% .
1	L	156	 % 51% 38% 9% .
1	M	156	 53% 37% 10% .
1	N	156	 53% 37% 9% .
1	O	156	 51% 40% 8% .
1	P	156	 53% 37% 8% .
1	Q	156	 % 53% 36% 10% .
1	R	156	 % 53% 37% 8% .
1	S	156	 54% 37% 8% .
1	T	156	 % 54% 36% 9% .
1	U	156	 % 54% 37% 8% .



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Mol	Chain	Length	Quality of chain
1	V	156	 54% 35% 10% .
1	W	156	 54% 37% 8% .
1	X	156	 56% 33% 9% .
1	Y	156	 2% 54% 37% 8% .
1	Z	156	 58% 33% 8% .
1	a	156	 1% 88% 10% .
1	b	156	 89% 10% .
1	c	156	 88% 11% .
1	d	156	 1% 87% 12% .
1	e	156	 1% 87% 12% .
1	f	156	 88% 11% .
1	g	156	 1% 87% 12% .
1	h	156	 89% 10% .
1	i	156	 88% 11% .
1	j	156	 88% 11% .
1	k	156	 87% 12% .
1	l	156	 88% 11% .
1	m	156	 1% 88% 11% .
1	n	156	 88% 11% .
1	o	156	 88% 11% .
1	p	156	 88% 10% .
1	q	156	 88% 10% .
1	r	156	 1% 88% 10% .
1	s	156	 88% 11% .
1	t	156	 88% 11% .

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Mol	Chain	Length	Quality of chain
1	u	156	 88% 11% .
1	v	156	 88% 10% .
1	w	156	 88% 11% .
1	x	156	 88% 10% .
1	y	156	 88% 11% .

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	SO4	5	157	-	-	X	-
2	SO4	6	157	-	-	X	-
2	SO4	7	158	-	-	X	-
2	SO4	B	158	-	-	X	-
2	SO4	B	159	-	-	X	-
2	SO4	H	158	-	-	X	-
2	SO4	L	158	-	-	-	X
2	SO4	P	157	-	-	X	-

## 2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 66544 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a protein called 6,7-dimethyl-8-ribityllumazine synthase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	154	Total	C	N	O	S	0	0	0
			1102	696	187	217	2			
1	B	154	Total	C	N	O	S	0	0	0
			1102	696	187	217	2			
1	C	154	Total	C	N	O	S	0	0	0
			1102	696	187	217	2			
1	D	154	Total	C	N	O	S	0	0	0
			1106	699	188	217	2			
1	E	154	Total	C	N	O	S	0	0	0
			1102	696	187	217	2			
1	F	154	Total	C	N	O	S	0	0	0
			1102	696	187	217	2			
1	G	154	Total	C	N	O	S	0	0	0
			1102	696	187	217	2			
1	H	154	Total	C	N	O	S	0	0	0
			1102	696	187	217	2			
1	I	154	Total	C	N	O	S	0	0	0
			1102	696	187	217	2			
1	J	154	Total	C	N	O	S	0	0	0
			1102	696	187	217	2			
1	K	154	Total	C	N	O	S	0	0	0
			1102	696	187	217	2			
1	L	154	Total	C	N	O	S	0	0	0
			1102	696	187	217	2			
1	M	154	Total	C	N	O	S	0	0	0
			1102	696	187	217	2			
1	N	154	Total	C	N	O	S	0	0	0
			1102	696	187	217	2			
1	O	154	Total	C	N	O	S	0	0	0
			1102	696	187	217	2			
1	P	154	Total	C	N	O	S	0	0	0
			1099	695	187	215	2			

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	Q	154	Total	C	N	O	S	0	0	0
			1102	696	187	217	2			
1	R	154	Total	C	N	O	S	0	0	0
			1102	696	187	217	2			
1	S	154	Total	C	N	O	S	0	0	0
			1102	696	187	217	2			
1	T	154	Total	C	N	O	S	0	0	0
			1102	696	187	217	2			
1	U	154	Total	C	N	O	S	0	0	0
			1102	696	187	217	2			
1	V	154	Total	C	N	O	S	0	0	0
			1102	696	187	217	2			
1	W	154	Total	C	N	O	S	0	0	0
			1102	696	187	217	2			
1	X	154	Total	C	N	O	S	0	0	0
			1102	696	187	217	2			
1	Y	154	Total	C	N	O	S	0	0	0
			1099	695	187	215	2			
1	Z	154	Total	C	N	O	S	0	0	0
			1102	696	187	217	2			
1	1	154	Total	C	N	O	S	0	0	0
			1106	699	188	217	2			
1	2	154	Total	C	N	O	S	0	0	0
			1102	696	187	217	2			
1	3	154	Total	C	N	O	S	0	0	0
			1102	696	187	217	2			
1	4	154	Total	C	N	O	S	0	0	0
			1102	696	187	217	2			
1	5	154	Total	C	N	O	S	0	0	0
			1106	699	188	217	2			
1	6	154	Total	C	N	O	S	0	0	0
			1106	699	188	217	2			
1	7	154	Total	C	N	O	S	0	0	0
			1102	696	187	217	2			
1	8	154	Total	C	N	O	S	0	0	0
			1102	696	187	217	2			
1	9	154	Total	C	N	O	S	0	0	0
			1106	699	188	217	2			
1	a	154	Total	C	N	O	S	0	0	0
			1102	696	187	217	2			
1	b	154	Total	C	N	O	S	0	0	0
			1106	699	188	217	2			

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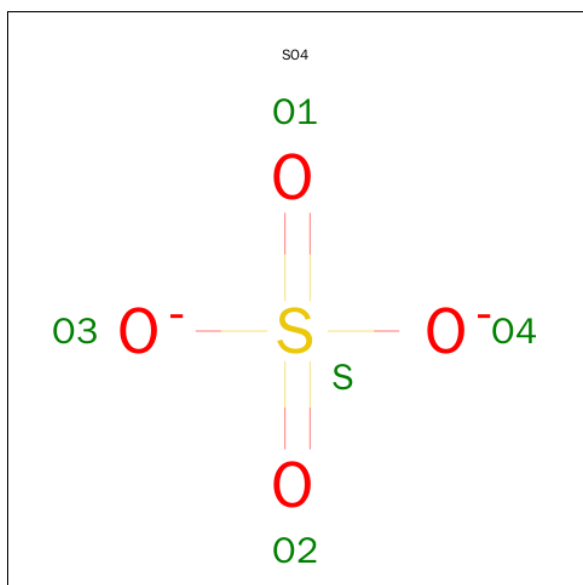
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	c	154	Total	C	N	O	S	0	0	0
			1102	696	187	217	2			
1	d	154	Total	C	N	O	S	0	0	0
			1102	696	187	217	2			
1	e	154	Total	C	N	O	S	0	0	0
			1102	696	187	217	2			
1	f	154	Total	C	N	O	S	0	0	0
			1106	699	188	217	2			
1	g	154	Total	C	N	O	S	0	0	0
			1106	699	188	217	2			
1	h	154	Total	C	N	O	S	0	0	0
			1102	696	187	217	2			
1	i	154	Total	C	N	O	S	0	0	0
			1102	696	187	217	2			
1	j	154	Total	C	N	O	S	0	0	0
			1106	699	188	217	2			
1	k	154	Total	C	N	O	S	0	0	0
			1102	696	187	217	2			
1	l	154	Total	C	N	O	S	0	0	0
			1106	699	188	217	2			
1	m	154	Total	C	N	O	S	0	0	0
			1102	696	187	217	2			
1	n	154	Total	C	N	O	S	0	0	0
			1102	696	187	217	2			
1	o	154	Total	C	N	O	S	0	0	0
			1099	695	187	215	2			
1	p	154	Total	C	N	O	S	0	0	0
			1102	696	187	217	2			
1	q	154	Total	C	N	O	S	0	0	0
			1102	696	187	217	2			
1	r	154	Total	C	N	O	S	0	0	0
			1102	696	187	217	2			
1	s	154	Total	C	N	O	S	0	0	0
			1102	696	187	217	2			
1	t	154	Total	C	N	O	S	0	0	0
			1102	696	187	217	2			
1	u	154	Total	C	N	O	S	0	0	0
			1099	695	187	215	2			
1	v	154	Total	C	N	O	S	0	0	0
			1102	696	187	217	2			
1	w	154	Total	C	N	O	S	0	0	0
			1102	696	187	217	2			

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	x	154	Total	C	N	O	S	0	0	0
			1099	695	187	215	2			
1	y	154	Total	C	N	O	S	0	0	0
			1106	699	188	217	2			

- Molecule 2 is SULFATE ION (three-letter code: SO4) (formula: O<sub>4</sub>S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	B	1	Total	O	S	0	0
			5	4	1		
2	B	1	Total	O	S	0	0
			5	4	1		
2	E	1	Total	O	S	0	0
			5	4	1		
2	A	1	Total	O	S	0	0
			5	4	1		
2	D	1	Total	O	S	0	0
			5	4	1		
2	G	1	Total	O	S	0	0
			5	4	1		
2	H	1	Total	O	S	0	0
			5	4	1		
2	I	1	Total	O	S	0	0
			5	4	1		
2	K	1	Total	O	S	0	0
			5	4	1		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	L	1	Total	O	S	0	0
			5	4	1		
2	M	1	Total	O	S	0	0
			5	4	1		
2	M	1	Total	O	S	0	0
			5	4	1		
2	O	1	Total	O	S	0	0
			5	4	1		
2	P	1	Total	O	S	0	0
			5	4	1		
2	P	1	Total	O	S	0	0
			5	4	1		
2	R	1	Total	O	S	0	0
			5	4	1		
2	R	1	Total	O	S	0	0
			5	4	1		
2	T	1	Total	O	S	0	0
			5	4	1		
2	U	1	Total	O	S	0	0
			5	4	1		
2	V	1	Total	O	S	0	0
			5	4	1		
2	W	1	Total	O	S	0	0
			5	4	1		
2	Z	1	Total	O	S	0	0
			5	4	1		
2	1	1	Total	O	S	0	0
			5	4	1		
2	1	1	Total	O	S	0	0
			5	4	1		
2	2	1	Total	O	S	0	0
			5	4	1		
2	3	1	Total	O	S	0	0
			5	4	1		
2	5	1	Total	O	S	0	0
			5	4	1		
2	6	1	Total	O	S	0	0
			5	4	1		
2	7	1	Total	O	S	0	0
			5	4	1		
2	Y	1	Total	O	S	0	0
			5	4	1		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	8	1	Total	O	S	0	0
			5	4	1		
2	9	1	Total	O	S	0	0
			5	4	1		
2	a	1	Total	O	S	0	0
			5	4	1		
2	b	1	Total	O	S	0	0
			5	4	1		
2	b	1	Total	O	S	0	0
			5	4	1		
2	d	1	Total	O	S	0	0
			5	4	1		
2	e	1	Total	O	S	0	0
			5	4	1		
2	f	1	Total	O	S	0	0
			5	4	1		
2	g	1	Total	O	S	0	0
			5	4	1		
2	h	1	Total	O	S	0	0
			5	4	1		
2	h	1	Total	O	S	0	0
			5	4	1		
2	i	1	Total	O	S	0	0
			5	4	1		
2	l	1	Total	O	S	0	0
			5	4	1		
2	m	1	Total	O	S	0	0
			5	4	1		
2	n	1	Total	O	S	0	0
			5	4	1		
2	o	1	Total	O	S	0	0
			5	4	1		
2	p	1	Total	O	S	0	0
			5	4	1		
2	p	1	Total	O	S	0	0
			5	4	1		
2	s	1	Total	O	S	0	0
			5	4	1		
2	s	1	Total	O	S	0	0
			5	4	1		
2	v	1	Total	O	S	0	0
			5	4	1		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	x	1	Total 5	O 4	S 1	0	0
2	x	1	Total 5	O 4	S 1	0	0
2	y	1	Total 5	O 4	S 1	0	0
2	F	1	Total 5	O 4	S 1	0	0
2	X	1	Total 5	O 4	S 1	0	0
2	k	1	Total 5	O 4	S 1	0	0
2	J	1	Total 5	O 4	S 1	0	0
2	o	1	Total 5	O 4	S 1	0	0
2	p	1	Total 5	O 4	S 1	0	0
2	q	1	Total 5	O 4	S 1	0	0
2	q	1	Total 5	O 4	S 1	0	0
2	Z	1	Total 5	O 4	S 1	0	0
2	B	1	Total 5	O 4	S 1	0	0
2	2	1	Total 5	O 4	S 1	0	0
2	h	1	Total 5	O 4	S 1	0	0
2	D	1	Total 5	O 4	S 1	0	0
2	F	1	Total 5	O 4	S 1	0	0
2	H	1	Total 5	O 4	S 1	0	0
2	L	1	Total 5	O 4	S 1	0	0
2	L	1	Total 5	O 4	S 1	0	0
2	M	1	Total 5	O 4	S 1	0	0

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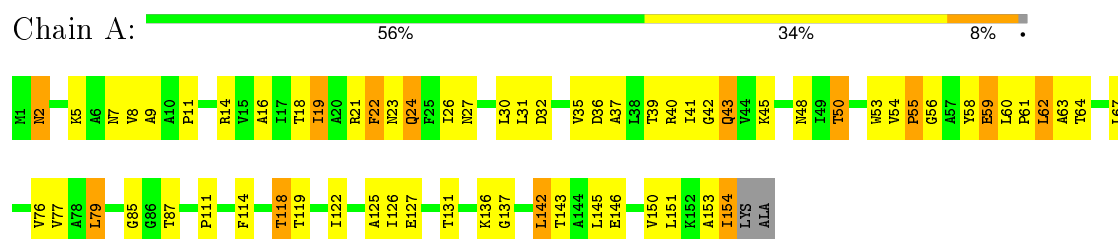
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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	Q	1	Total	O	S	0	0
			5	4	1		
2	7	1	Total	O	S	0	0
			5	4	1		
2	b	1	Total	O	S	0	0
			5	4	1		
2	d	1	Total	O	S	0	0
			5	4	1		
2	k	1	Total	O	S	0	0
			5	4	1		
2	u	1	Total	O	S	0	0
			5	4	1		
2	1	1	Total	O	S	0	0
			5	4	1		

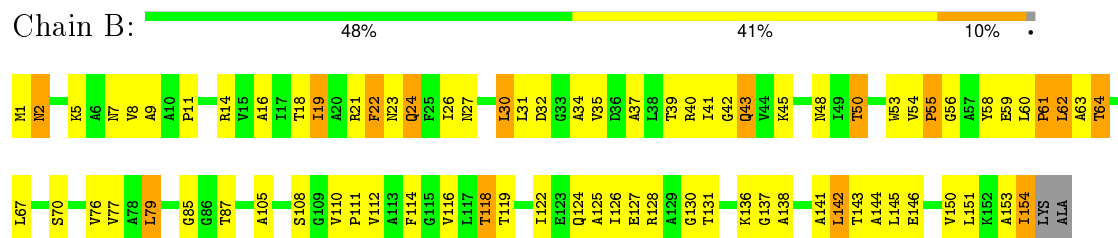
### 3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ( $RSRZ > 2$ ). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

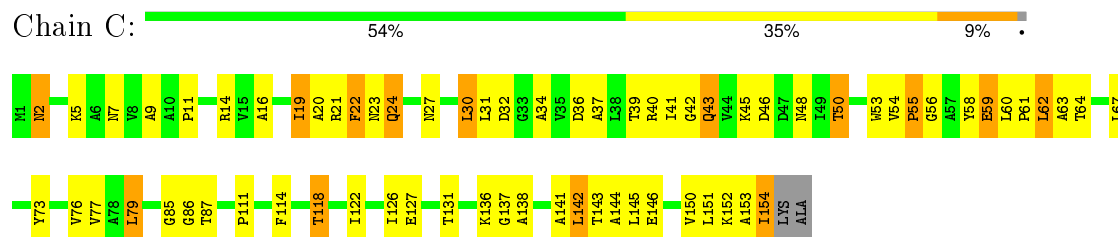
- Molecule 1: 6,7-dimethyl-8-ribityllumazine synthase



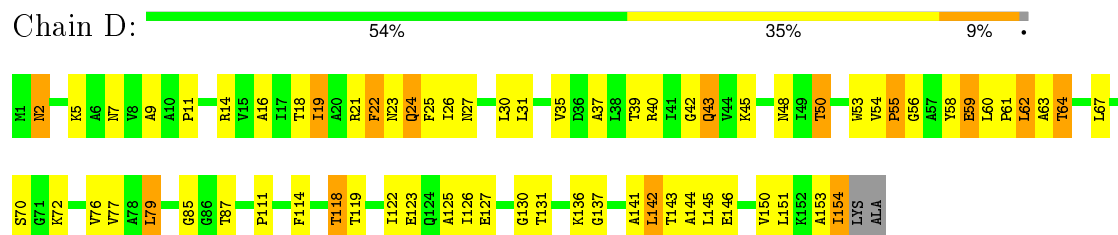
- Molecule 1: 6,7-dimethyl-8-ribityllumazine synthase



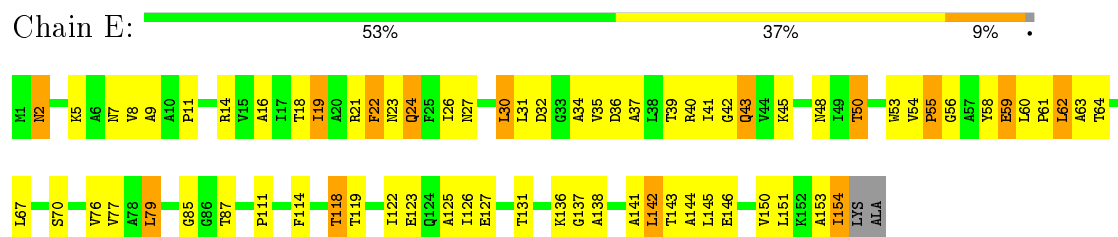
- Molecule 1: 6,7-dimethyl-8-ribityllumazine synthase



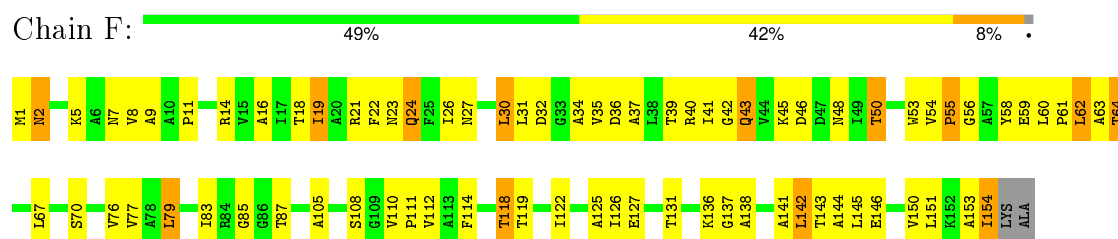
- Molecule 1: 6,7-dimethyl-8-ribityllumazine synthase



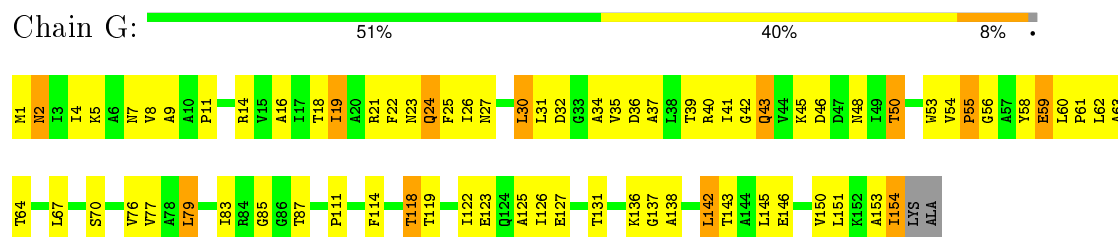
- Molecule 1: 6,7-dimethyl-8-ribityllumazine synthase



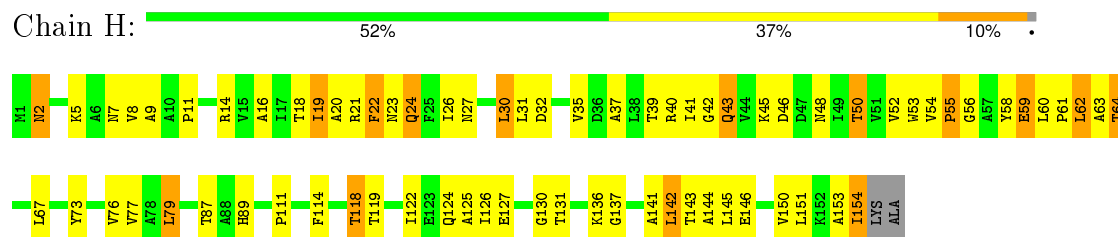
- Molecule 1: 6,7-dimethyl-8-ribityllumazine synthase



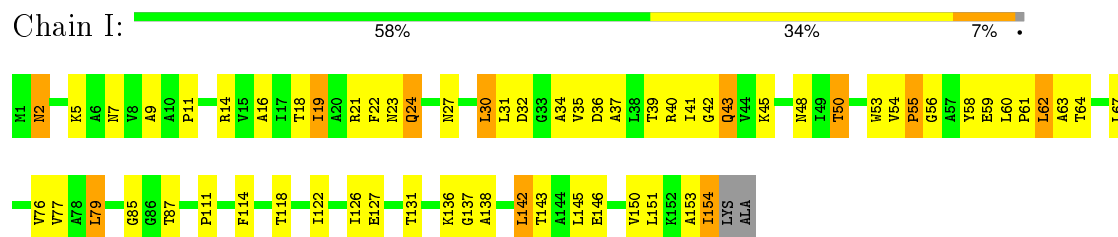
- Molecule 1: 6,7-dimethyl-8-ribityllumazine synthase



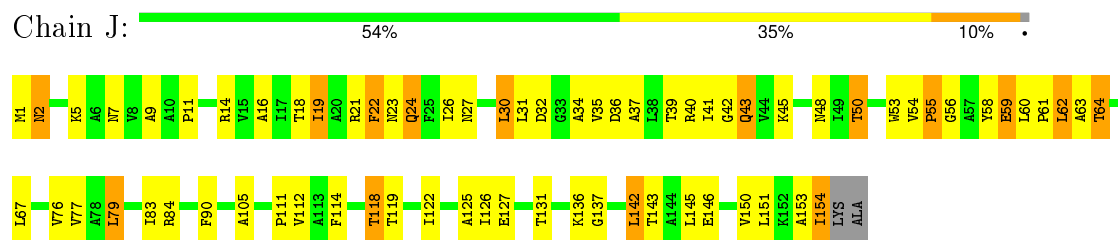
- Molecule 1: 6,7-dimethyl-8-ribityllumazine synthase



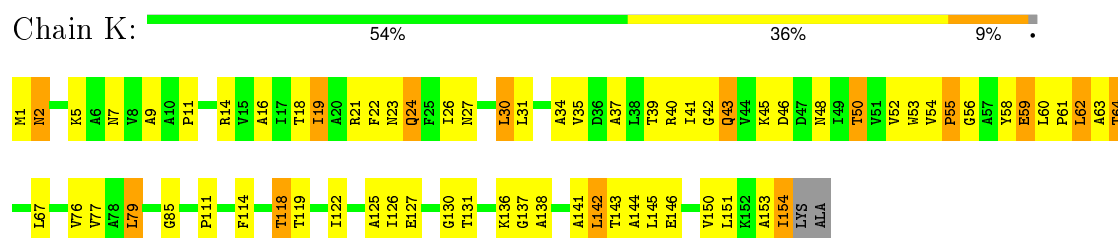
- Molecule 1: 6,7-dimethyl-8-ribityllumazine synthase



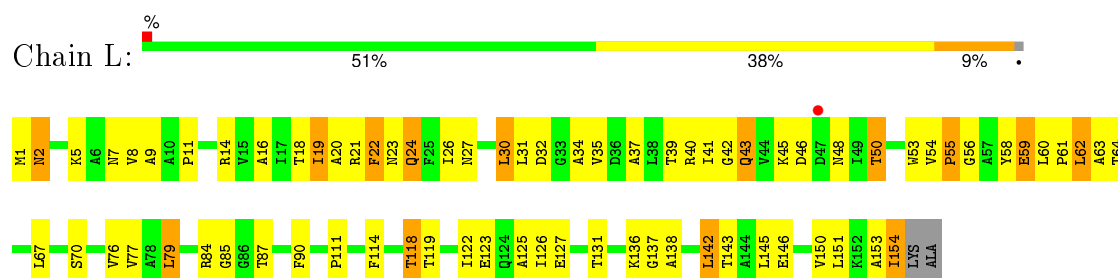
- Molecule 1: 6,7-dimethyl-8-ribityllumazine synthase



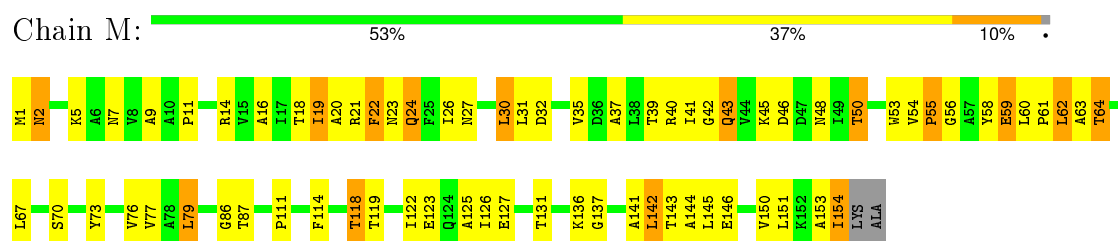
- Molecule 1: 6,7-dimethyl-8-ribityllumazine synthase



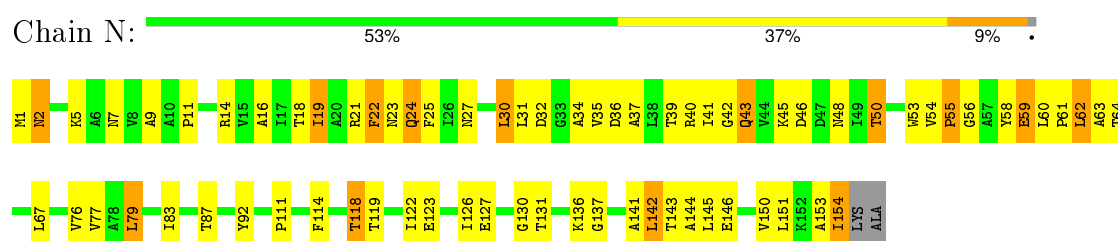
- Molecule 1: 6,7-dimethyl-8-ribityllumazine synthase



- Molecule 1: 6,7-dimethyl-8-ribityllumazine synthase

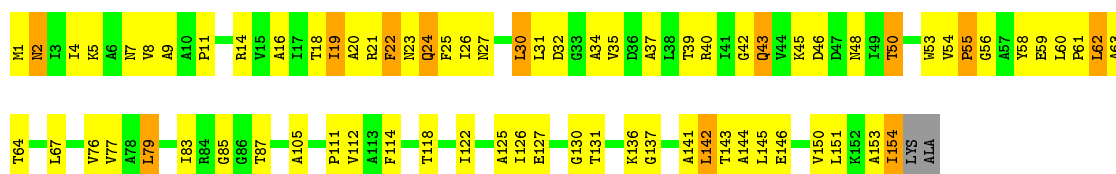


- Molecule 1: 6,7-dimethyl-8-ribityllumazine synthase



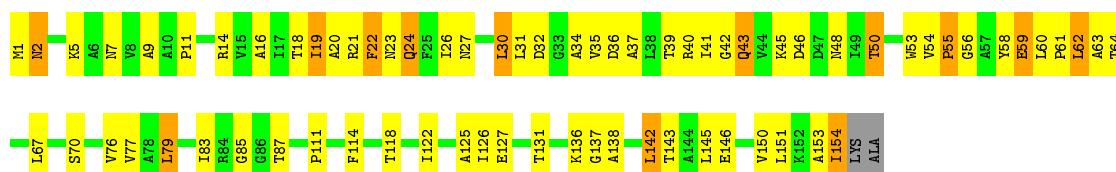
- Molecule 1: 6,7-dimethyl-8-ribityllumazine synthase





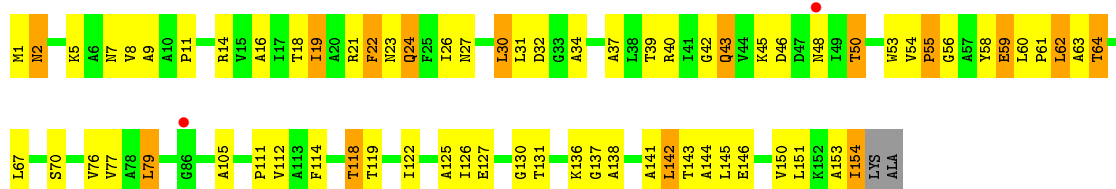
- Molecule 1: 6,7-dimethyl-8-ribityllumazine synthase

Chain P: 53% 37% 8% .



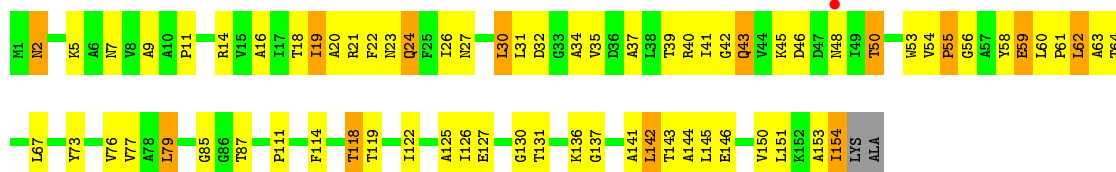
- Molecule 1: 6,7-dimethyl-8-ribityllumazine synthase

Chain Q: 53% 36% 10% .



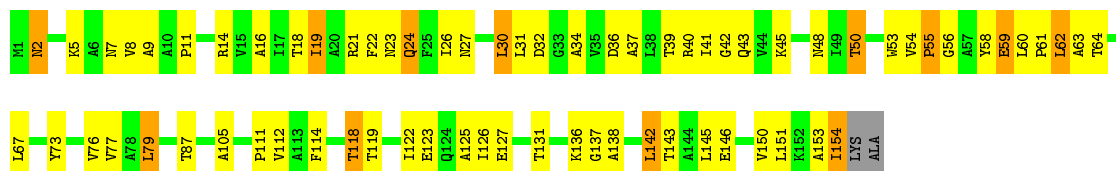
- Molecule 1: 6,7-dimethyl-8-ribityllumazine synthase

Chain R: 53% 37% 8% .



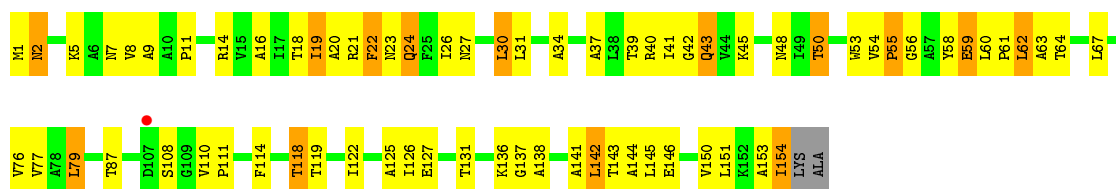
- Molecule 1: 6,7-dimethyl-8-ribityllumazine synthase

Chain S: 54% 37% 8% .

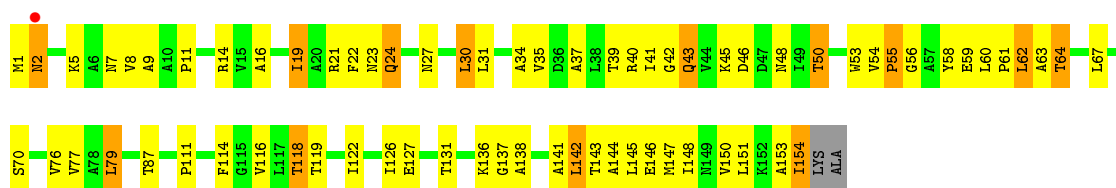


- Molecule 1: 6,7-dimethyl-8-ribityllumazine synthase

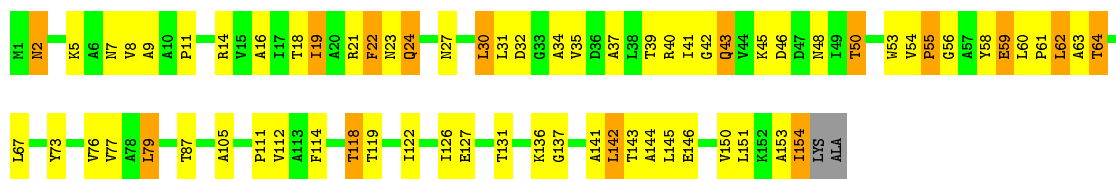
Chain T: 54% 36% 9% .



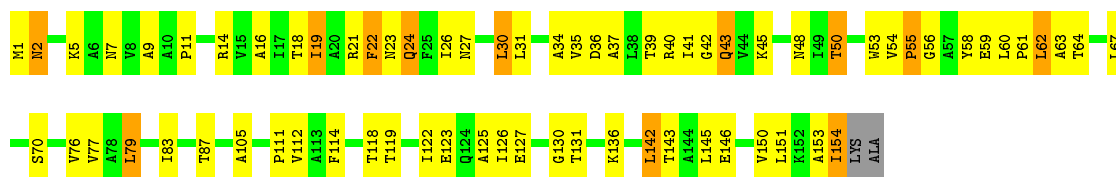
- Molecule 1: 6,7-dimethyl-8-ribityllumazine synthase



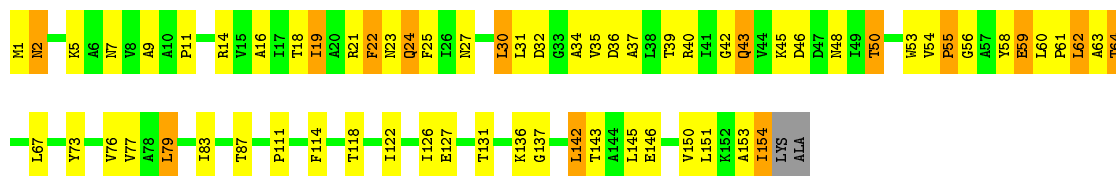
- Molecule 1: 6,7-dimethyl-8-ribityllumazine synthase



- Molecule 1: 6,7-dimethyl-8-ribityllumazine synthase

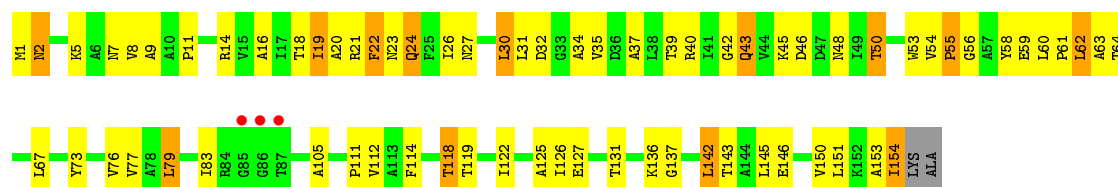


- Molecule 1: 6,7-dimethyl-8-ribityllumazine synthase



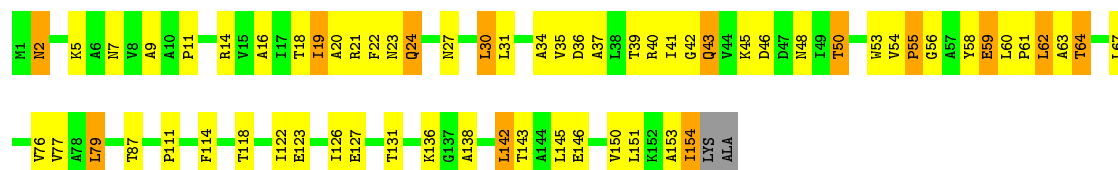
- Molecule 1: 6,7-dimethyl-8-ribityllumazine synthase





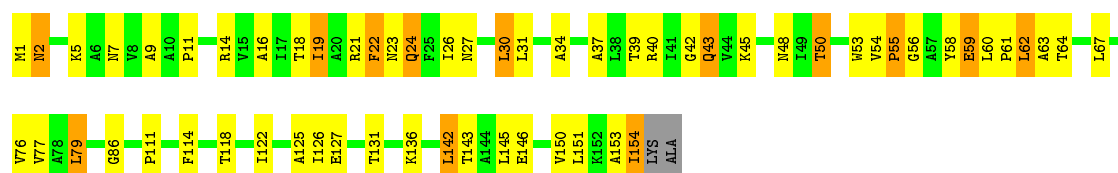
- Molecule 1: 6,7-dimethyl-8-ribityllumazine synthase

Chain Z: 58% 33% 8%



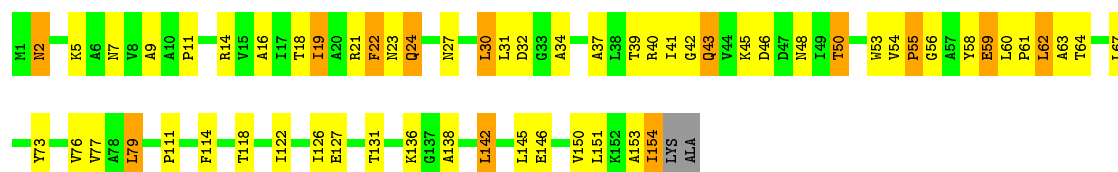
- Molecule 1: 6,7-dimethyl-8-ribityllumazine synthase

Chain 1: 60% 30% 8%



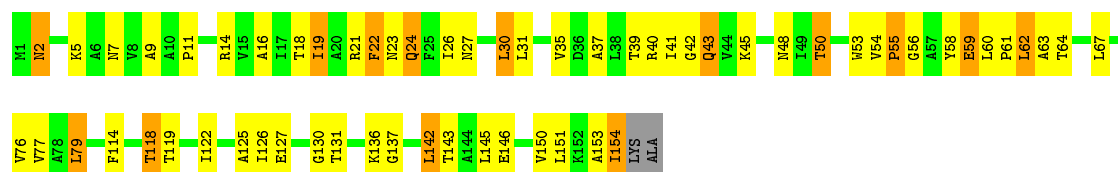
- Molecule 1: 6,7-dimethyl-8-ribityllumazine synthase

Chain 2: 60% 30% 8%



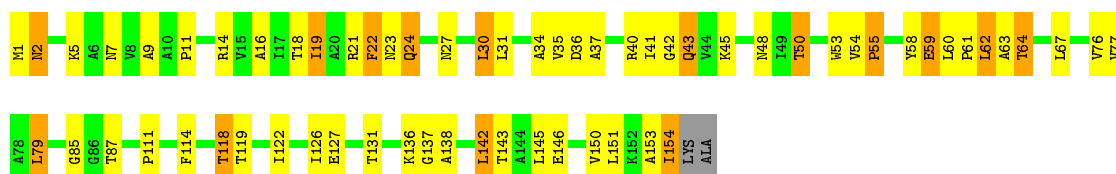
- Molecule 1: 6,7-dimethyl-8-ribityllumazine synthase

Chain 3: 60% 30% 9%



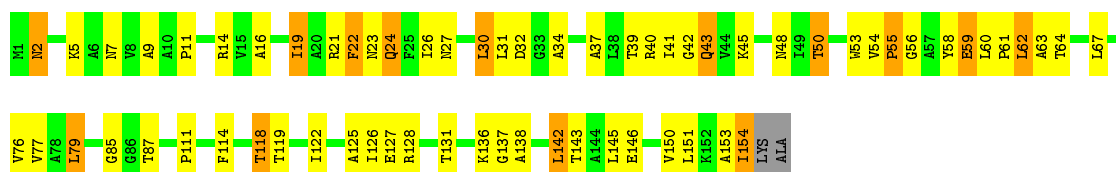
- Molecule 1: 6,7-dimethyl-8-ribityllumazine synthase

Chain 4: 58% 31% 10%



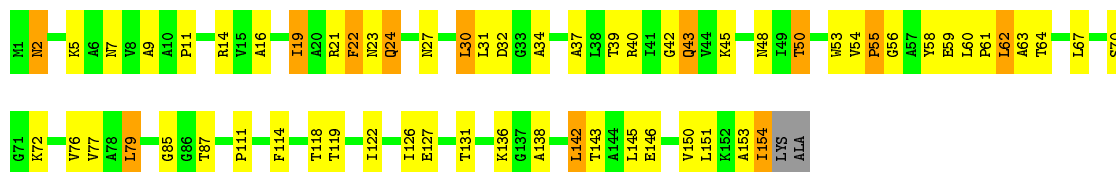
- Molecule 1: 6,7-dimethyl-8-ribityllumazine synthase

Chain 5: 57% 33% 9% .



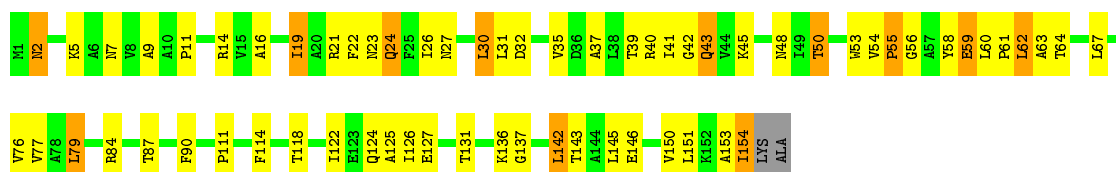
- Molecule 1: 6,7-dimethyl-8-ribityllumazine synthase

Chain 6: 59% 32% 8% .



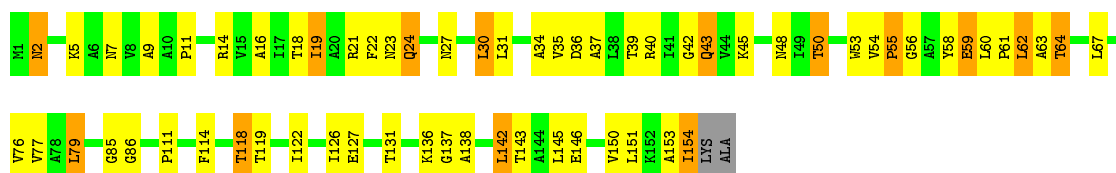
- Molecule 1: 6,7-dimethyl-8-ribityllumazine synthase

Chain 7: 58% 33% 8% .



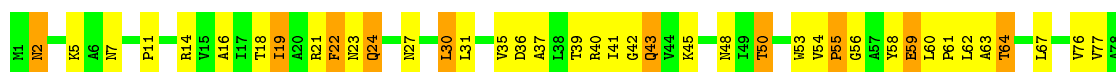
- Molecule 1: 6,7-dimethyl-8-ribityllumazine synthase

Chain 8: 58% 31% 9% .

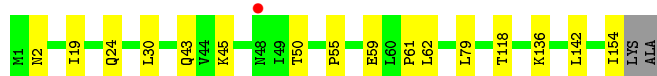
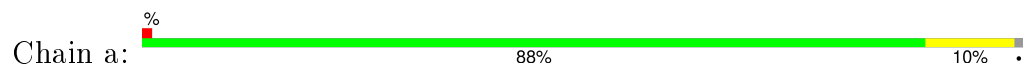


- Molecule 1: 6,7-dimethyl-8-ribityllumazine synthase

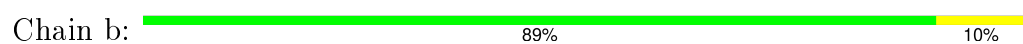
Chain 9: 60% 30% 9% .



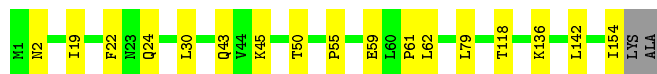
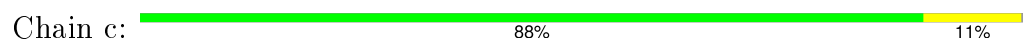
- Molecule 1: 6,7-dimethyl-8-ribityllumazine synthase



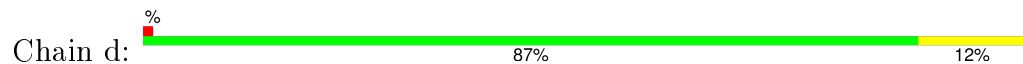
- Molecule 1: 6,7-dimethyl-8-ribityllumazine synthase



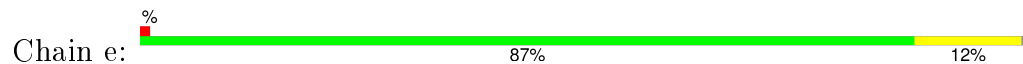
- Molecule 1: 6,7-dimethyl-8-ribityllumazine synthase



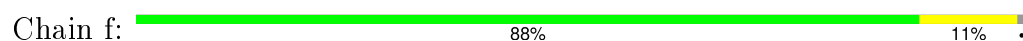
- Molecule 1: 6,7-dimethyl-8-ribityllumazine synthase



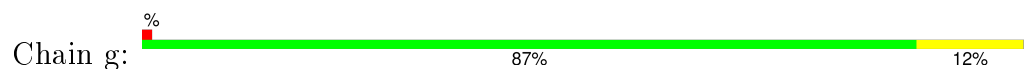
- Molecule 1: 6,7-dimethyl-8-ribityllumazine synthase



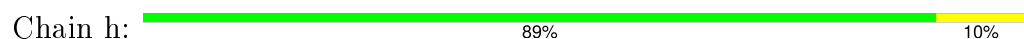
- Molecule 1: 6,7-dimethyl-8-ribityllumazine synthase



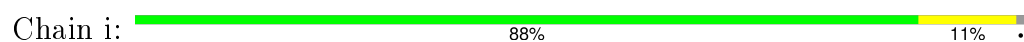
- Molecule 1: 6,7-dimethyl-8-ribityllumazine synthase



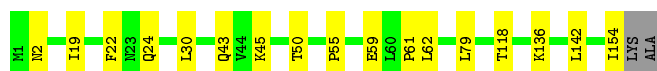
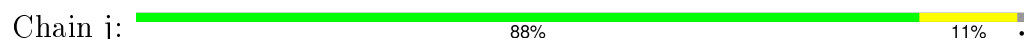
- Molecule 1: 6,7-dimethyl-8-ribityllumazine synthase



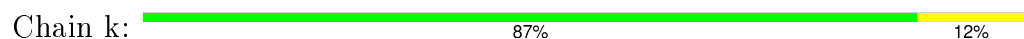
- Molecule 1: 6,7-dimethyl-8-ribityllumazine synthase



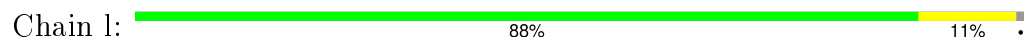
- Molecule 1: 6,7-dimethyl-8-ribityllumazine synthase



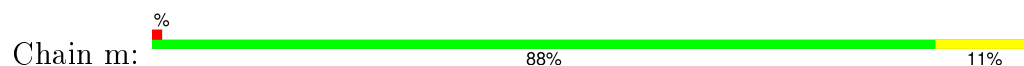
- Molecule 1: 6,7-dimethyl-8-ribityllumazine synthase



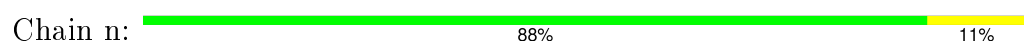
- Molecule 1: 6,7-dimethyl-8-ribityllumazine synthase



- Molecule 1: 6,7-dimethyl-8-ribityllumazine synthase



- Molecule 1: 6,7-dimethyl-8-ribityllumazine synthase





- Molecule 1: 6,7-dimethyl-8-ribityllumazine synthase

Chain o: 88% 11%



- Molecule 1: 6,7-dimethyl-8-ribityllumazine synthase

Chain p: 88% 10%



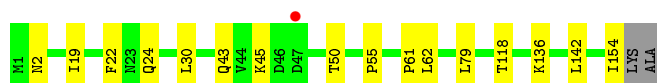
- Molecule 1: 6,7-dimethyl-8-ribityllumazine synthase

Chain q: 88% 10%



- Molecule 1: 6,7-dimethyl-8-ribityllumazine synthase

Chain r: 88% 10%



- Molecule 1: 6,7-dimethyl-8-ribityllumazine synthase

Chain s: 88% 11%



- Molecule 1: 6,7-dimethyl-8-ribityllumazine synthase

Chain t: 88% 11%



- Molecule 1: 6,7-dimethyl-8-ribityllumazine synthase

Chain u: 88% 11%



- Molecule 1: 6,7-dimethyl-8-ribityllumazine synthase

Chain v: 88% 10% .



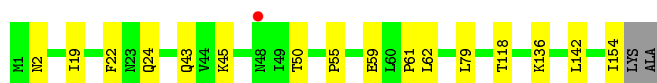
- Molecule 1: 6,7-dimethyl-8-ribityllumazine synthase

Chain w: 88% 11% .



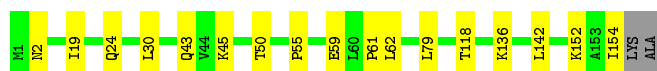
- Molecule 1: 6,7-dimethyl-8-ribityllumazine synthase

Chain x: 88% 10% .



- Molecule 1: 6,7-dimethyl-8-ribityllumazine synthase

Chain y: 88% 11% .



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	154.23Å 151.50Å 235.03Å 90.00° 97.08° 90.00°	Depositor
Resolution (Å)	32.03 – 3.57 32.03 – 3.57	Depositor EDS
% Data completeness (in resolution range)	82.5 (32.03-3.57) 82.5 (32.03-3.57)	Depositor EDS
$R_{merge}$	0.09	Depositor
$R_{sym}$	0.09	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.58 (at 3.56Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.6 _289)	Depositor
R, $R_{free}$	0.226 , 0.264 0.208 , 0.247	Depositor DCC
$R_{free}$ test set	1052 reflections (1.00%)	DCC
Wilson B-factor (Å <sup>2</sup> )	87.6	Xtriage
Anisotropy	0.413	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.30 , 44.6	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	2 of 105298 reflections (0.002%)	Xtriage
$F_o, F_c$ correlation	0.91	EDS
Total number of atoms	66544	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	81.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.81% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical values of  $\langle |L| \rangle$ ,  $\langle L^2 \rangle$  for acentric reflections are 0.5, 0.375 respectively for untwinned datasets, and 0.333, 0.2 for perfectly twinned datasets.

## 5 Model quality

### 5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: SO4

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with  $|Z| > 5$  is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z  > 5$	RMSZ	$\# Z  > 5$
1	1	0.46	0/1119	0.57	0/1524
1	2	0.45	0/1115	0.56	0/1520
1	3	0.43	0/1115	0.54	0/1520
1	4	0.42	0/1115	0.54	0/1520
1	5	0.48	0/1119	0.56	0/1524
1	6	0.42	0/1119	0.55	0/1524
1	7	0.45	0/1115	0.57	0/1520
1	8	0.44	0/1115	0.57	0/1520
1	9	0.46	0/1119	0.55	0/1524
1	A	0.45	0/1115	0.56	0/1520
1	B	0.45	0/1115	0.56	0/1520
1	C	0.43	0/1115	0.58	0/1520
1	D	0.46	0/1119	0.57	0/1524
1	E	0.45	0/1115	0.57	0/1520
1	F	0.43	0/1115	0.56	0/1520
1	G	0.42	0/1115	0.56	0/1520
1	H	0.41	0/1115	0.56	0/1520
1	I	0.42	0/1115	0.55	0/1520
1	J	0.43	0/1115	0.54	0/1520
1	K	0.44	0/1115	0.56	0/1520
1	L	0.43	0/1115	0.56	0/1520
1	M	0.41	0/1115	0.54	0/1520
1	N	0.46	0/1115	0.56	0/1520
1	O	0.47	0/1115	0.55	0/1520
1	P	0.41	0/1112	0.56	0/1516
1	Q	0.44	0/1115	0.55	0/1520
1	R	0.44	0/1115	0.56	0/1520
1	S	0.44	0/1115	0.54	0/1520
1	T	0.44	0/1115	0.56	0/1520
1	U	0.44	0/1115	0.56	0/1520
1	V	0.43	0/1115	0.56	0/1520
1	W	0.42	0/1115	0.54	0/1520

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z  >5	RMSZ	# Z  >5
1	X	0.44	0/1115	0.55	0/1520
1	Y	0.43	0/1112	0.54	0/1516
1	Z	0.42	0/1115	0.55	0/1520
1	a	0.43	0/1115	0.55	0/1520
1	b	0.45	0/1119	0.56	0/1524
1	c	0.42	0/1115	0.55	0/1520
1	d	0.41	0/1115	0.54	0/1520
1	e	0.42	0/1115	0.55	0/1520
1	f	0.42	0/1119	0.55	0/1524
1	g	0.44	0/1119	0.56	0/1524
1	h	0.42	0/1115	0.55	0/1520
1	i	0.43	0/1115	0.55	0/1520
1	j	0.43	0/1119	0.57	0/1524
1	k	0.43	0/1115	0.57	0/1520
1	l	0.44	0/1119	0.56	0/1524
1	m	0.44	0/1115	0.56	0/1520
1	n	0.41	0/1115	0.56	0/1520
1	o	0.42	0/1112	0.55	0/1516
1	p	0.44	0/1115	0.55	0/1520
1	q	0.44	0/1115	0.55	0/1520
1	r	0.44	0/1115	0.56	0/1520
1	s	0.43	0/1115	0.56	0/1520
1	t	0.42	0/1115	0.57	0/1520
1	u	0.42	0/1112	0.55	0/1516
1	v	0.41	0/1115	0.53	0/1520
1	w	0.43	0/1115	0.54	0/1520
1	x	0.39	0/1112	0.54	0/1516
1	y	0.43	0/1119	0.56	0/1524
All	All	0.43	0/66929	0.55	0/91224

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

## 5.2 Too-close contacts ⓘ

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	1	1106	0	1126	56	0
1	2	1102	0	1115	58	0
1	3	1102	0	1115	59	0
1	4	1102	0	1115	60	0
1	5	1106	0	1126	59	0
1	6	1106	0	1126	53	0
1	7	1102	0	1115	54	0
1	8	1102	0	1115	61	0
1	9	1106	0	1126	59	0
1	A	1102	0	1115	72	0
1	B	1102	0	1115	80	0
1	C	1102	0	1115	75	0
1	D	1106	0	1126	72	0
1	E	1102	0	1115	76	0
1	F	1102	0	1115	85	0
1	G	1102	0	1115	82	0
1	H	1102	0	1115	79	0
1	I	1102	0	1115	74	0
1	J	1102	0	1115	76	0
1	K	1102	0	1115	75	0
1	L	1102	0	1115	86	0
1	M	1102	0	1115	92	0
1	N	1102	0	1115	83	0
1	O	1102	0	1115	83	0
1	P	1099	0	1113	89	0
1	Q	1102	0	1115	76	0
1	R	1102	0	1115	86	0
1	S	1102	0	1115	76	0
1	T	1102	0	1115	81	0
1	U	1102	0	1115	83	0
1	V	1102	0	1115	81	0
1	W	1102	0	1115	83	0
1	X	1102	0	1115	76	0
1	Y	1099	0	1113	72	0
1	Z	1102	0	1115	65	0
1	a	1102	0	1115	0	0
1	b	1106	0	1126	0	0
1	c	1102	0	1115	0	0
1	d	1102	0	1115	0	0
1	e	1102	0	1115	0	0
1	f	1106	0	1126	0	0
1	g	1106	0	1126	0	0
1	h	1102	0	1115	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	i	1102	0	1115	0	0
1	j	1106	0	1126	0	0
1	k	1102	0	1115	0	0
1	l	1106	0	1126	0	0
1	m	1102	0	1115	0	0
1	n	1102	0	1115	0	0
1	o	1099	0	1113	0	0
1	p	1102	0	1115	0	0
1	q	1102	0	1115	0	0
1	r	1102	0	1115	0	0
1	s	1102	0	1115	0	0
1	t	1102	0	1115	0	0
1	u	1099	0	1113	0	0
1	v	1102	0	1115	0	0
1	w	1102	0	1115	0	0
1	x	1099	0	1113	0	0
1	y	1106	0	1126	0	0
2	1	15	0	0	0	0
2	2	10	0	0	0	0
2	3	5	0	0	1	0
2	5	5	0	0	2	0
2	6	5	0	0	2	0
2	7	10	0	0	2	0
2	8	5	0	0	1	0
2	9	5	0	0	0	0
2	A	5	0	0	1	0
2	B	15	0	0	6	0
2	D	10	0	0	1	0
2	E	5	0	0	1	0
2	F	10	0	0	1	0
2	G	5	0	0	1	0
2	H	10	0	0	3	0
2	I	5	0	0	1	0
2	J	5	0	0	0	0
2	K	5	0	0	1	0
2	L	15	0	0	2	0
2	M	15	0	0	0	0
2	O	5	0	0	1	0
2	P	10	0	0	2	0
2	Q	5	0	0	1	0
2	R	10	0	0	1	0
2	T	5	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	U	5	0	0	0	0
2	V	5	0	0	0	0
2	W	5	0	0	0	0
2	X	5	0	0	0	0
2	Y	5	0	0	0	0
2	Z	10	0	0	1	0
2	a	5	0	0	0	0
2	b	15	0	0	0	0
2	d	10	0	0	0	0
2	e	5	0	0	0	0
2	f	5	0	0	0	0
2	g	5	0	0	0	0
2	h	15	0	0	0	0
2	i	5	0	0	0	0
2	k	10	0	0	0	0
2	l	5	0	0	0	0
2	m	5	0	0	0	0
2	n	5	0	0	0	0
2	o	10	0	0	0	0
2	p	15	0	0	0	0
2	q	10	0	0	0	0
2	s	10	0	0	0	0
2	u	5	0	0	0	0
2	v	5	0	0	0	0
2	x	10	0	0	0	0
2	y	5	0	0	0	0
All	All	66544	0	67011	2088	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 16.

All (2088) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:Z:60:LEU:O	1:Z:64:THR:HG23	1.60	1.01
1:I:5:LYS:HE2	1:J:21:ARG:HH22	1.49	1.00
1:I:60:LEU:O	1:I:64:THR:HG23	1.79	0.98
1:G:60:LEU:O	1:G:64:THR:HG23	1.69	0.98
1:D:60:LEU:O	1:D:64:THR:HG23	1.64	0.97
1:6:122:ILE:H	1:6:122:ILE:HD12	1.30	0.96
1:U:60:LEU:O	1:U:64:THR:HG23	1.73	0.95

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:J:60:LEU:O	1:J:64:THR:HG23	1.70	0.94
1:M:122:ILE:HD12	1:M:122:ILE:H	1.33	0.94
1:X:60:LEU:O	1:X:64:THR:HG23	1.66	0.94
1:9:60:LEU:O	1:9:64:THR:HG23	1.68	0.94
1:5:60:LEU:O	1:5:64:THR:HG23	1.67	0.94
1:N:60:LEU:O	1:N:64:THR:HG23	1.70	0.94
1:E:60:LEU:O	1:E:64:THR:HG23	1.76	0.94
1:H:60:LEU:O	1:H:64:THR:HG23	1.67	0.93
1:1:60:LEU:O	1:1:64:THR:HG23	1.69	0.93
1:F:60:LEU:O	1:F:64:THR:HG23	1.77	0.93
1:P:60:LEU:O	1:P:64:THR:HG23	1.68	0.93
1:D:122:ILE:HD12	1:D:122:ILE:H	1.34	0.93
1:Z:122:ILE:HD12	1:Z:122:ILE:H	1.32	0.93
1:5:122:ILE:HD12	1:5:122:ILE:H	1.34	0.93
1:B:122:ILE:HD12	1:B:122:ILE:H	1.39	0.92
1:G:122:ILE:H	1:G:122:ILE:HD12	1.37	0.92
1:A:60:LEU:O	1:A:64:THR:HG23	1.71	0.92
1:7:60:LEU:O	1:7:64:THR:HG23	1.69	0.92
1:3:122:ILE:H	1:3:122:ILE:HD12	1.34	0.92
1:O:60:LEU:O	1:O:64:THR:HG23	1.76	0.92
1:K:122:ILE:H	1:K:122:ILE:HD12	1.37	0.92
1:2:60:LEU:O	1:2:64:THR:HG23	1.70	0.92
1:L:122:ILE:H	1:L:122:ILE:HD12	1.35	0.92
1:V:122:ILE:HD12	1:V:122:ILE:H	1.35	0.91
1:Q:122:ILE:HD12	1:Q:122:ILE:H	1.41	0.91
1:R:60:LEU:O	1:R:64:THR:HG23	1.70	0.91
1:M:60:LEU:O	1:M:64:THR:HG23	1.71	0.91
1:C:5:LYS:HE2	1:D:21:ARG:HH22	1.36	0.91
1:V:5:LYS:HE2	1:W:21:ARG:HH22	1.34	0.91
1:I:122:ILE:H	1:I:122:ILE:HD12	1.36	0.90
1:W:122:ILE:HD12	1:W:122:ILE:H	1.38	0.90
1:9:122:ILE:HD12	1:9:122:ILE:H	1.36	0.90
1:S:60:LEU:O	1:S:64:THR:HG23	1.76	0.90
1:J:122:ILE:HD12	1:J:122:ILE:H	1.36	0.90
1:H:122:ILE:HD12	1:H:122:ILE:H	1.36	0.90
1:6:60:LEU:O	1:6:64:THR:HG23	1.71	0.90
1:4:60:LEU:O	1:4:64:THR:HG23	1.71	0.89
1:R:122:ILE:HD12	1:R:122:ILE:H	1.38	0.89
1:V:60:LEU:O	1:V:64:THR:HG23	1.76	0.89
1:7:122:ILE:HD12	1:7:122:ILE:H	1.35	0.89
1:S:122:ILE:H	1:S:122:ILE:HD12	1.38	0.89

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:O:122:ILE:HD12	1:O:122:ILE:H	1.38	0.89
1:T:122:ILE:HD12	1:T:122:ILE:H	1.41	0.88
1:Y:60:LEU:O	1:Y:64:THR:HG23	1.73	0.88
1:B:60:LEU:O	1:B:64:THR:HG23	1.73	0.88
1:Y:122:ILE:H	1:Y:122:ILE:HD12	1.38	0.88
1:Q:60:LEU:O	1:Q:64:THR:HG23	1.73	0.88
1:A:122:ILE:HD12	1:A:122:ILE:H	1.39	0.88
1:F:122:ILE:H	1:F:122:ILE:HD12	1.41	0.88
1:E:122:ILE:HD12	1:E:122:ILE:H	1.40	0.88
1:U:5:LYS:HE2	1:V:21:ARG:HH22	1.56	0.88
1:N:122:ILE:HD12	1:N:122:ILE:H	1.40	0.88
1:1:122:ILE:HD12	1:1:122:ILE:H	1.38	0.88
1:R:5:LYS:HE2	1:S:21:ARG:HH22	1.39	0.88
1:8:60:LEU:O	1:8:64:THR:HG23	1.72	0.88
1:2:122:ILE:HD12	1:2:122:ILE:H	1.38	0.88
1:U:122:ILE:HD12	1:U:122:ILE:H	1.39	0.87
1:T:60:LEU:O	1:T:64:THR:HG23	1.77	0.87
1:C:60:LEU:O	1:C:64:THR:HG23	1.73	0.87
1:X:122:ILE:HD12	1:X:122:ILE:H	1.38	0.87
1:P:122:ILE:HD12	1:P:122:ILE:H	1.38	0.87
2:B:158:SO4:O2	1:C:86:GLY:HA3	1.97	0.87
1:W:60:LEU:O	1:W:64:THR:HG23	1.74	0.87
1:C:122:ILE:HD12	1:C:122:ILE:H	1.40	0.86
1:A:21:ARG:HH22	1:E:5:LYS:HE2	1.44	0.86
1:5:5:LYS:HE2	1:6:21:ARG:HH22	1.39	0.86
1:8:122:ILE:H	1:8:122:ILE:HD12	1.38	0.86
1:D:5:LYS:HE2	1:E:21:ARG:HH22	1.41	0.86
1:L:60:LEU:O	1:L:64:THR:HG23	1.75	0.85
1:3:60:LEU:O	1:3:64:THR:HG23	1.77	0.85
1:F:21:ARG:HH22	1:J:5:LYS:HE2	1.43	0.84
1:4:122:ILE:H	1:4:122:ILE:HD12	1.41	0.83
1:S:5:LYS:HE2	1:T:21:ARG:HH22	1.42	0.83
1:B:5:LYS:HE2	1:C:21:ARG:HH22	1.41	0.83
1:6:85:GLY:HA3	2:6:157:SO4:O1	1.76	0.83
1:2:5:LYS:HE2	1:3:21:ARG:HH22	1.44	0.82
2:Z:158:SO4:O2	1:1:86:GLY:HA3	1.81	0.81
1:K:60:LEU:O	1:K:64:THR:HG23	1.80	0.81
1:7:5:LYS:HE2	1:8:21:ARG:HH22	1.46	0.81
1:G:5:LYS:HE2	1:H:21:ARG:HH22	1.44	0.80
1:Z:5:LYS:HE2	1:1:21:ARG:HH22	1.45	0.80
1:L:5:LYS:HE2	1:M:21:ARG:HH22	1.43	0.80

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:8:5:LYS:HE2	1:9:21:ARG:HH22	1.45	0.80
1:A:5:LYS:HE2	1:B:21:ARG:HH22	1.47	0.80
1:P:5:LYS:HE2	1:Q:21:ARG:HH22	1.45	0.79
1:Q:5:LYS:HE2	1:R:21:ARG:HH22	1.48	0.79
1:U:50:THR:HG22	1:Y:2:ASN:HD21	1.48	0.79
1:5:21:ARG:HH22	1:9:5:LYS:HE2	1.46	0.79
1:M:5:LYS:HE2	1:N:21:ARG:HH22	1.71	0.79
1:P:85:GLY:HA3	2:P:157:SO4:O4	1.84	0.77
1:K:50:THR:HG22	1:O:2:ASN:HD21	1.67	0.77
1:W:5:LYS:HE2	1:X:21:ARG:HH22	1.49	0.77
1:H:5:LYS:HE2	1:I:21:ARG:HH22	1.58	0.76
1:D:85:GLY:HA3	2:D:157:SO4:O2	2.87	0.75
1:I:85:GLY:HA3	2:I:157:SO4:O1	1.87	0.75
1:F:114:PHE:HB2	1:G:58:TYR:CE2	2.39	0.75
1:B:22:PHE:O	1:B:23:ASN:HB2	1.96	0.75
1:U:21:ARG:HH22	1:Y:5:LYS:HE2	1.50	0.75
1:C:146:GLU:O	1:C:150:VAL:HG23	1.87	0.74
1:M:114:PHE:HB2	1:N:58:TYR:CE2	2.22	0.74
1:U:114:PHE:HB2	1:V:58:TYR:CE2	2.22	0.74
1:K:21:ARG:HH22	1:O:5:LYS:HE2	1.53	0.74
1:6:5:LYS:HE2	1:7:21:ARG:HH22	1.52	0.74
1:G:23:ASN:HA	1:G:131:THR:HA	17.50	0.74
1:M:131:THR:HA	1:X:23:ASN:HA	78.28	0.73
1:G:2:ASN:HD21	1:H:50:THR:HG22	1.54	0.73
1:2:22:PHE:O	1:2:23:ASN:HB2	1.88	0.73
1:K:122:ILE:O	1:K:126:ILE:HG12	1.95	0.72
1:A:2:ASN:HD21	1:B:50:THR:HG22	1.85	0.72
1:E:122:ILE:O	1:E:126:ILE:HG12	1.89	0.72
1:M:146:GLU:O	1:M:150:VAL:HG23	1.90	0.72
1:X:114:PHE:HB2	1:Y:58:TYR:CE2	2.37	0.72
1:I:122:ILE:O	1:I:126:ILE:HG12	2.00	0.71
1:3:5:LYS:HE2	1:4:21:ARG:HH22	1.53	0.71
1:V:22:PHE:O	1:V:23:ASN:HB2	1.93	0.71
1:F:5:LYS:HE2	1:G:21:ARG:HH22	1.56	0.71
1:O:22:PHE:O	1:O:23:ASN:HB2	1.90	0.71
1:K:58:TYR:CE2	1:O:114:PHE:HB2	2.29	0.71
1:F:58:TYR:CE2	1:J:114:PHE:HB2	2.36	0.71
1:W:122:ILE:O	1:W:126:ILE:HG12	1.98	0.71
1:Z:21:ARG:HH22	1:4:5:LYS:HE2	1.55	0.70
1:V:146:GLU:O	1:V:150:VAL:HG23	1.91	0.70
1:X:5:LYS:HE2	1:Y:21:ARG:HH22	1.58	0.70

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:N:23:ASN:HA	1:W:131:THR:HA	70.61	0.70
1:K:146:GLU:O	1:K:150:VAL:HG23	1.96	0.70
1:4:122:ILE:O	1:4:126:ILE:HG12	1.92	0.70
1:K:22:PHE:O	1:K:23:ASN:HB2	1.91	0.70
1:F:146:GLU:O	1:F:150:VAL:HG23	2.06	0.70
1:L:2:ASN:HD21	1:M:50:THR:HG22	1.69	0.69
1:5:22:PHE:O	1:5:23:ASN:HB2	1.93	0.69
1:4:146:GLU:O	1:4:150:VAL:HG23	1.93	0.69
1:R:122:ILE:O	1:R:126:ILE:HG12	2.05	0.69
1:K:114:PHE:HB2	1:L:58:TYR:CE2	2.40	0.69
1:N:2:ASN:HD21	1:O:50:THR:HG22	1.58	0.69
1:3:22:PHE:O	1:3:23:ASN:HB2	1.92	0.69
1:H:122:ILE:O	1:H:126:ILE:HG12	1.93	0.69
1:P:21:ARG:HH22	1:T:5:LYS:HE2	1.64	0.68
1:M:22:PHE:O	1:M:23:ASN:HB2	2.01	0.68
1:D:122:ILE:O	1:D:126:ILE:HG12	1.93	0.68
1:5:85:GLY:HA3	2:5:157:SO4:O2	1.93	0.68
1:G:122:ILE:O	1:G:126:ILE:HG12	1.98	0.68
1:N:5:LYS:HE2	1:O:21:ARG:HH22	1.59	0.68
1:P:122:ILE:O	1:P:126:ILE:HG12	1.96	0.68
1:L:122:ILE:O	1:L:126:ILE:HG12	2.06	0.68
1:S:114:PHE:HB2	1:T:58:TYR:CE2	2.36	0.68
1:V:2:ASN:HD21	1:W:50:THR:HG22	1.59	0.68
1:E:146:GLU:O	1:E:150:VAL:HG23	1.93	0.68
1:G:22:PHE:O	1:G:23:ASN:HB2	1.98	0.68
1:G:154:ILE:HD13	1:H:67:LEU:HD21	2.03	0.68
1:A:122:ILE:O	1:A:126:ILE:HG12	1.94	0.68
1:T:146:GLU:O	1:T:150:VAL:HG23	1.94	0.68
1:1:114:PHE:HB2	1:2:58:TYR:CE2	2.29	0.68
1:U:119:THR:HG22	1:V:87:THR:HG21	1.74	0.67
1:P:146:GLU:O	1:P:150:VAL:HG23	1.93	0.67
1:Q:122:ILE:O	1:Q:126:ILE:HG12	1.94	0.67
1:1:14:ARG:HA	1:1:48:ASN:HB3	1.76	0.67
1:3:146:GLU:O	1:3:150:VAL:HG23	1.94	0.67
1:Z:122:ILE:O	1:Z:126:ILE:HG12	1.94	0.67
1:P:58:TYR:CE2	1:T:114:PHE:HB2	2.30	0.67
1:F:22:PHE:O	1:F:23:ASN:HB2	1.94	0.67
1:I:22:PHE:O	1:I:23:ASN:HB2	1.94	0.67
1:R:146:GLU:O	1:R:150:VAL:HG23	1.95	0.67
1:S:122:ILE:O	1:S:126:ILE:HG12	2.00	0.67
1:O:122:ILE:O	1:O:126:ILE:HG12	1.95	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:14:ARG:HA	1:A:48:ASN:HB3	1.77	0.67
1:P:50:THR:HG22	1:T:2:ASN:HD21	1.60	0.67
1:H:114:PHE:HB2	1:I:58:TYR:CE2	2.30	0.67
1:I:5:LYS:HE2	1:J:21:ARG:NH2	2.25	0.66
1:T:122:ILE:O	1:T:126:ILE:HG12	2.03	0.66
1:P:114:PHE:HB2	1:Q:58:TYR:CE2	2.40	0.66
1:X:22:PHE:O	1:X:23:ASN:HB2	1.94	0.66
1:J:122:ILE:O	1:J:126:ILE:HG12	1.94	0.66
1:F:122:ILE:O	1:F:126:ILE:HG12	1.99	0.66
1:5:122:ILE:O	1:5:126:ILE:HG12	1.95	0.66
1:1:122:ILE:O	1:1:126:ILE:HG12	1.96	0.66
1:Y:146:GLU:O	1:Y:150:VAL:HG23	2.05	0.66
1:8:122:ILE:O	1:8:126:ILE:HG12	1.96	0.66
1:W:1:MET:HB2	1:X:46:ASP:O	2.41	0.66
1:8:146:GLU:O	1:8:150:VAL:HG23	1.96	0.66
1:7:22:PHE:O	1:7:23:ASN:HB2	1.94	0.66
1:4:22:PHE:O	1:4:23:ASN:HB2	1.96	0.66
1:J:14:ARG:HA	1:J:48:ASN:HB3	1.80	0.66
1:U:122:ILE:O	1:U:126:ILE:HG12	1.95	0.66
1:P:2:ASN:HD21	1:Q:50:THR:HG22	1.65	0.66
1:5:62:LEU:HD13	1:9:151:LEU:HD11	1.78	0.66
1:2:122:ILE:O	1:2:126:ILE:HG12	1.94	0.66
1:G:151:LEU:HD11	1:H:62:LEU:HD13	2.02	0.66
1:L:22:PHE:O	1:L:23:ASN:HB2	1.96	0.66
1:K:5:LYS:HE2	1:L:21:ARG:HH22	1.61	0.66
1:X:146:GLU:O	1:X:150:VAL:HG23	1.94	0.66
1:W:22:PHE:O	1:W:23:ASN:HB2	1.98	0.66
1:L:14:ARG:HA	1:L:48:ASN:HB3	1.78	0.66
1:H:124:GLN:HG2	2:H:158:SO4:O2	2.46	0.65
1:W:114:PHE:HB2	1:X:58:TYR:CE2	2.35	0.65
1:U:146:GLU:O	1:U:150:VAL:HG23	1.96	0.65
1:J:146:GLU:O	1:J:150:VAL:HG23	2.00	0.65
1:P:46:ASP:O	1:T:1:MET:HB2	1.96	0.65
1:3:122:ILE:O	1:3:126:ILE:HG12	1.95	0.65
1:A:50:THR:HG22	1:E:2:ASN:HD21	1.61	0.65
1:T:14:ARG:HA	1:T:48:ASN:HB3	1.78	0.65
1:N:114:PHE:HB2	1:O:58:TYR:CE2	2.38	0.65
1:L:146:GLU:O	1:L:150:VAL:HG23	1.97	0.65
1:I:2:ASN:HD21	1:J:50:THR:HG22	1.61	0.65
1:X:14:ARG:HA	1:X:48:ASN:HB3	1.78	0.65
1:B:146:GLU:O	1:B:150:VAL:HG23	2.00	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:37:ALA:HB1	1:E:142:LEU:HD13	1.78	0.65
1:V:122:ILE:O	1:V:126:ILE:HG12	1.96	0.65
1:W:14:ARG:HA	1:W:48:ASN:HB3	1.79	0.65
1:Q:114:PHE:HB2	1:R:58:TYR:CE2	2.32	0.65
1:U:14:ARG:HA	1:U:48:ASN:HB3	1.79	0.65
1:I:5:LYS:CE	1:J:21:ARG:HH22	2.30	0.65
1:E:131:THR:HA	1:O:23:ASN:HA	1.78	0.65
1:B:14:ARG:HA	1:B:48:ASN:HB3	1.79	0.65
1:Z:58:TYR:CE2	1:4:114:PHE:HB2	2.31	0.65
1:8:14:ARG:HA	1:8:48:ASN:HB3	1.78	0.65
1:Q:7:ASN:ND2	1:Z:41:ILE:HA	118.75	0.65
1:I:146:GLU:O	1:I:150:VAL:HG23	1.97	0.65
1:M:14:ARG:HA	1:M:48:ASN:HB3	1.78	0.65
1:C:22:PHE:O	1:C:23:ASN:HB2	2.00	0.65
1:V:114:PHE:HB2	1:W:58:TYR:CE2	2.41	0.65
1:Q:146:GLU:O	1:Q:150:VAL:HG23	1.97	0.65
1:G:114:PHE:HB2	1:H:58:TYR:CE2	2.52	0.65
1:E:14:ARG:HA	1:E:48:ASN:HB3	1.78	0.65
1:C:2:ASN:HD21	1:D:50:THR:HG22	1.77	0.65
1:G:1:MET:HA	1:H:46:ASP:O	2.20	0.64
1:X:122:ILE:O	1:X:126:ILE:HG12	2.01	0.64
1:F:14:ARG:HA	1:F:48:ASN:HB3	1.79	0.64
1:S:22:PHE:O	1:S:23:ASN:HB2	2.04	0.64
1:V:5:LYS:HE2	1:W:21:ARG:NH2	2.10	0.64
1:R:14:ARG:HA	1:R:48:ASN:HB3	1.83	0.64
1:P:55:PRO:HB2	1:T:143:THR:HG22	1.79	0.64
1:D:146:GLU:O	1:D:150:VAL:HG23	2.02	0.64
1:Y:122:ILE:O	1:Y:126:ILE:HG12	1.98	0.64
1:J:22:PHE:O	1:J:23:ASN:HB2	1.97	0.64
2:Q:158:SO4:O4	1:R:85:GLY:HA3	9.45	0.64
1:6:122:ILE:H	1:6:122:ILE:CD1	2.07	0.64
1:D:22:PHE:O	1:D:23:ASN:HB2	2.01	0.64
1:J:37:ALA:HB1	1:J:142:LEU:HD13	1.85	0.64
1:G:14:ARG:HA	1:G:48:ASN:HB3	1.80	0.64
1:6:22:PHE:O	1:6:23:ASN:HB2	1.98	0.64
1:N:14:ARG:HA	1:N:48:ASN:HB3	1.79	0.64
1:N:122:ILE:O	1:N:126:ILE:HG12	1.98	0.64
1:K:143:THR:HG22	1:L:55:PRO:HB2	1.96	0.64
1:W:146:GLU:O	1:W:150:VAL:HG23	1.97	0.64
1:8:22:PHE:O	1:8:23:ASN:HB2	1.96	0.64
1:6:122:ILE:O	1:6:126:ILE:HG12	1.98	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:K:14:ARG:HA	1:K:48:ASN:HB3	1.83	0.64
1:6:14:ARG:HA	1:6:48:ASN:HB3	1.80	0.64
1:D:14:ARG:HA	1:D:48:ASN:HB3	1.80	0.64
1:4:14:ARG:HA	1:4:48:ASN:HB3	1.80	0.64
1:N:146:GLU:O	1:N:150:VAL:HG23	1.98	0.64
1:A:22:PHE:O	1:A:23:ASN:HB2	1.97	0.64
1:W:2:ASN:HD21	1:X:50:THR:HG22	1.77	0.64
1:C:14:ARG:HA	1:C:48:ASN:HB3	1.80	0.64
1:7:122:ILE:O	1:7:126:ILE:HG12	1.98	0.63
1:T:22:PHE:O	1:T:23:ASN:HB2	1.97	0.63
1:U:58:TYR:CE2	1:Y:114:PHE:HB2	2.35	0.63
1:C:114:PHE:HB2	1:D:58:TYR:CE2	2.45	0.63
1:6:146:GLU:O	1:6:150:VAL:HG23	1.98	0.63
1:H:14:ARG:HA	1:H:48:ASN:HB3	1.80	0.63
1:U:22:PHE:O	1:U:23:ASN:HB2	1.98	0.63
1:3:122:ILE:CD1	1:3:122:ILE:H	2.10	0.63
1:Y:14:ARG:HA	1:Y:48:ASN:HB3	1.80	0.63
1:P:22:PHE:O	1:P:23:ASN:HB2	1.99	0.63
1:O:14:ARG:HA	1:O:48:ASN:HB3	1.81	0.63
1:V:143:THR:HG22	1:W:55:PRO:HB2	2.06	0.63
1:Z:22:PHE:O	1:Z:23:ASN:HB2	1.98	0.63
1:L:122:ILE:H	1:L:122:ILE:CD1	2.10	0.63
1:G:37:ALA:HB1	1:G:142:LEU:HD13	1.81	0.63
1:7:114:PHE:HB2	1:8:58:TYR:CE2	2.34	0.63
1:F:37:ALA:HB1	1:F:142:LEU:HD13	1.81	0.63
1:C:122:ILE:O	1:C:126:ILE:HG12	1.99	0.63
1:Z:146:GLU:O	1:Z:150:VAL:HG23	1.99	0.63
1:O:146:GLU:O	1:O:150:VAL:HG23	2.04	0.63
1:S:2:ASN:HD21	1:T:50:THR:HG22	1.63	0.63
1:1:22:PHE:O	1:1:23:ASN:HB2	1.98	0.62
1:U:118:THR:O	1:V:87:THR:HB	1.99	0.62
1:2:11:PRO:HB3	1:2:43:GLN:HB3	1.81	0.62
1:H:146:GLU:O	1:H:150:VAL:HG23	1.98	0.62
2:L:158:SO4:O4	1:M:86:GLY:HA3	1.98	0.62
1:5:87:THR:HB	1:9:118:THR:O	1.99	0.62
1:R:11:PRO:HB3	1:R:43:GLN:HB3	1.81	0.62
1:7:14:ARG:HA	1:7:48:ASN:HB3	1.81	0.62
1:L:37:ALA:HB1	1:L:142:LEU:HD13	1.82	0.62
1:R:24:GLN:HA	1:R:27:ASN:HB2	1.81	0.62
1:N:22:PHE:O	1:N:23:ASN:HB2	2.00	0.62
1:S:146:GLU:O	1:S:150:VAL:HG23	2.05	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:Q:22:PHE:O	1:Q:23:ASN:HB2	2.01	0.62
1:K:122:ILE:H	1:K:122:ILE:CD1	2.13	0.62
1:T:122:ILE:H	1:T:122:ILE:CD1	2.17	0.62
1:R:22:PHE:O	1:R:23:ASN:HB2	2.02	0.62
1:9:122:ILE:O	1:9:126:ILE:HG12	1.99	0.62
1:A:146:GLU:O	1:A:150:VAL:HG23	1.99	0.62
1:Q:14:ARG:HA	1:Q:48:ASN:HB3	1.82	0.62
1:S:14:ARG:HA	1:S:48:ASN:HB3	1.80	0.61
1:V:14:ARG:HA	1:V:48:ASN:HB3	1.82	0.61
1:M:122:ILE:CD1	1:M:122:ILE:H	2.09	0.61
1:R:127:GLU:HA	1:R:131:THR:OG1	2.01	0.61
1:D:122:ILE:H	1:D:122:ILE:CD1	2.09	0.61
1:U:5:LYS:HE2	1:V:21:ARG:NH2	2.35	0.61
1:O:11:PRO:HB3	1:O:43:GLN:HB3	1.87	0.61
1:A:85:GLY:HA3	2:A:157:SO4:O1	1.99	0.61
1:M:122:ILE:O	1:M:126:ILE:HG12	1.99	0.61
1:B:122:ILE:O	1:B:126:ILE:HG12	2.02	0.61
1:E:77:VAL:HG12	1:E:79:LEU:HD13	1.84	0.61
1:7:24:GLN:HA	1:7:27:ASN:HB2	1.82	0.61
1:T:153:ALA:O	1:T:154:ILE:HG12	2.12	0.61
1:2:14:ARG:HA	1:2:48:ASN:HB3	1.81	0.61
1:M:37:ALA:HB1	1:M:142:LEU:HD13	1.83	0.61
1:Z:14:ARG:HA	1:Z:48:ASN:HB3	1.83	0.61
1:K:37:ALA:HB1	1:K:142:LEU:HD13	1.84	0.61
1:9:22:PHE:O	1:9:23:ASN:HB2	1.99	0.61
1:3:118:THR:O	1:4:87:THR:HB	2.01	0.61
1:K:118:THR:O	1:L:87:THR:HB	2.00	0.61
1:G:16:ALA:HB1	1:G:67:LEU:HD13	1.89	0.61
1:U:2:ASN:HD21	1:V:50:THR:HG22	1.71	0.61
1:5:2:ASN:HD21	1:6:50:THR:HG22	1.66	0.61
1:U:37:ALA:HB1	1:U:142:LEU:HD13	1.83	0.61
1:2:146:GLU:O	1:2:150:VAL:HG23	2.01	0.61
1:R:122:ILE:CD1	1:R:122:ILE:H	2.14	0.60
1:B:124:GLN:HG2	2:B:158:SO4:O4	2.01	0.60
1:K:53:TRP:CD1	1:O:5:LYS:HE3	2.56	0.60
1:G:146:GLU:O	1:G:150:VAL:HG23	2.01	0.60
1:H:22:PHE:O	1:H:23:ASN:HB2	2.02	0.60
1:L:114:PHE:HB2	1:M:58:TYR:CE2	2.36	0.60
1:5:14:ARG:HA	1:5:48:ASN:HB3	1.81	0.60
1:U:11:PRO:HB3	1:U:43:GLN:HB3	1.84	0.60
1:Q:127:GLU:HA	1:Q:131:THR:OG1	2.02	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:122:ILE:CD1	1:H:122:ILE:H	2.12	0.60
1:T:16:ALA:HB1	1:T:67:LEU:HD13	1.89	0.60
1:N:77:VAL:HG12	1:N:79:LEU:HD13	1.93	0.60
1:W:42:GLY:HA3	1:W:145:LEU:HD13	1.83	0.60
1:T:37:ALA:HB1	1:T:142:LEU:HD13	1.82	0.60
1:J:122:ILE:CD1	1:J:122:ILE:H	2.13	0.60
1:D:123:GLU:OE2	1:F:122:ILE:HD13	2.35	0.60
1:D:131:THR:HA	1:F:23:ASN:HA	2.14	0.60
1:D:114:PHE:HB2	1:E:58:TYR:CE2	2.35	0.60
1:4:127:GLU:HA	1:4:131:THR:OG1	2.02	0.60
1:5:122:ILE:CD1	1:5:122:ILE:H	2.10	0.60
1:N:153:ALA:O	1:N:154:ILE:HG12	2.10	0.60
1:P:87:THR:HG21	1:T:119:THR:HG22	1.84	0.60
1:B:37:ALA:HB1	1:B:142:LEU:HD13	1.86	0.60
1:9:14:ARG:HA	1:9:48:ASN:HB3	1.82	0.60
1:F:11:PRO:HB3	1:F:43:GLN:HB3	1.83	0.60
1:C:37:ALA:HB1	1:C:142:LEU:HD13	1.84	0.60
1:5:58:TYR:CE2	1:9:114:PHE:HB2	2.37	0.60
1:I:14:ARG:HA	1:I:48:ASN:HB3	1.85	0.60
1:R:23:ASN:HA	1:Z:131:THR:HA	88.27	0.60
1:E:22:PHE:O	1:E:23:ASN:HB2	2.01	0.60
1:Q:37:ALA:HB1	1:Q:142:LEU:HD13	1.84	0.60
1:D:11:PRO:HB3	1:D:43:GLN:HB3	1.91	0.59
1:I:11:PRO:HB3	1:I:43:GLN:HB3	1.84	0.59
1:S:11:PRO:HB3	1:S:43:GLN:HB3	1.83	0.59
1:N:37:ALA:HB1	1:N:142:LEU:HD13	1.84	0.59
1:6:37:ALA:HB1	1:6:142:LEU:HD13	1.83	0.59
1:S:127:GLU:HA	1:S:131:THR:OG1	2.02	0.59
1:Z:50:THR:HG22	1:4:2:ASN:HD21	1.67	0.59
1:H:2:ASN:HD21	1:I:50:THR:HG22	1.84	0.59
1:A:122:ILE:CD1	1:A:122:ILE:H	2.15	0.59
1:1:146:GLU:O	1:1:150:VAL:HG23	2.02	0.59
1:Q:118:THR:O	1:R:87:THR:HB	2.02	0.59
1:X:37:ALA:HB1	1:X:142:LEU:HD13	1.86	0.59
1:U:143:THR:HG22	1:V:55:PRO:HB2	1.83	0.59
1:H:118:THR:O	1:I:87:THR:HB	2.02	0.59
1:F:2:ASN:HD21	1:G:50:THR:HG22	1.76	0.59
1:1:5:LYS:HE2	1:2:21:ARG:HH22	1.68	0.59
1:O:37:ALA:HB1	1:O:142:LEU:HD13	1.89	0.59
1:D:37:ALA:HB1	1:D:142:LEU:HD13	1.85	0.59
1:V:37:ALA:HB1	1:V:142:LEU:HD13	1.87	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:V:122:ILE:H	1:V:122:ILE:CD1	2.11	0.59
1:2:122:ILE:CD1	1:2:122:ILE:H	2.14	0.59
1:A:37:ALA:HB1	1:A:142:LEU:HD13	1.85	0.59
1:Q:122:ILE:CD1	1:Q:122:ILE:H	2.16	0.59
1:O:77:VAL:HG12	1:O:79:LEU:HD13	1.92	0.59
1:3:11:PRO:HB3	1:3:43:GLN:HB3	1.84	0.59
1:A:127:GLU:HA	1:A:131:THR:OG1	2.09	0.59
1:A:2:ASN:H	1:A:2:ASN:HD22	1.51	0.58
1:6:114:PHE:HB2	1:7:58:TYR:CE2	2.37	0.58
1:H:41:ILE:HA	1:2:7:ASN:ND2	96.49	0.58
1:3:37:ALA:HB1	1:3:142:LEU:HD13	1.85	0.58
1:P:14:ARG:HA	1:P:48:ASN:HB3	1.84	0.58
1:K:56:GLY:O	1:K:59:GLU:HG2	2.03	0.58
1:3:2:ASN:HD21	1:4:50:THR:HG22	1.65	0.58
1:I:37:ALA:HB1	1:I:142:LEU:HD13	1.85	0.58
1:Q:24:GLN:HA	1:Q:27:ASN:HB2	1.87	0.58
1:D:127:GLU:HA	1:D:131:THR:OG1	2.04	0.58
1:Y:22:PHE:O	1:Y:23:ASN:HB2	2.03	0.58
1:U:154:ILE:HD13	1:V:67:LEU:HD21	2.05	0.58
1:P:37:ALA:HB1	1:P:142:LEU:HD13	1.86	0.58
1:F:127:GLU:HA	1:F:131:THR:OG1	2.07	0.58
1:U:127:GLU:HA	1:U:131:THR:OG1	2.04	0.58
1:E:42:GLY:HA3	1:E:145:LEU:HD13	1.96	0.58
1:I:41:ILE:HA	1:9:7:ASN:ND2	112.29	0.58
1:Y:11:PRO:HB3	1:Y:43:GLN:HB3	1.88	0.58
1:1:11:PRO:HB3	1:1:43:GLN:HB3	1.85	0.58
1:W:122:ILE:CD1	1:W:122:ILE:H	2.14	0.58
1:A:114:PHE:HB2	1:B:58:TYR:CE2	2.38	0.58
1:S:122:ILE:H	1:S:122:ILE:CD1	2.13	0.58
1:3:14:ARG:HA	1:3:48:ASN:HB3	1.84	0.58
1:L:122:ILE:N	1:L:122:ILE:HD12	2.12	0.58
1:O:122:ILE:CD1	1:O:122:ILE:H	2.13	0.58
1:R:77:VAL:HG12	1:R:79:LEU:HD13	1.85	0.58
1:H:37:ALA:HB1	1:H:142:LEU:HD13	1.86	0.58
1:6:11:PRO:HB3	1:6:43:GLN:HB3	1.85	0.58
1:W:77:VAL:HG12	1:W:79:LEU:HD13	1.88	0.58
1:R:114:PHE:HB2	1:S:58:TYR:CE2	2.52	0.58
1:5:127:GLU:HA	1:5:131:THR:OG1	2.04	0.58
1:5:146:GLU:O	1:5:150:VAL:HG23	2.03	0.58
1:7:11:PRO:HB3	1:7:43:GLN:HB3	1.86	0.58
1:H:127:GLU:HA	1:H:131:THR:OG1	2.04	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:16:ALA:HB1	1:H:67:LEU:HD13	1.95	0.58
1:S:16:ALA:HB1	1:S:67:LEU:HD13	1.86	0.58
1:T:7:ASN:ND2	1:3:40:ARG:O	2.37	0.57
1:J:127:GLU:HA	1:J:131:THR:OG1	2.03	0.57
1:K:11:PRO:HB3	1:K:43:GLN:HB3	1.86	0.57
1:V:11:PRO:HB3	1:V:43:GLN:HB3	1.90	0.57
1:V:54:VAL:HB	1:V:55:PRO:HD2	1.87	0.57
1:1:127:GLU:HA	1:1:131:THR:OG1	2.04	0.57
1:2:37:ALA:HB1	1:2:142:LEU:HD13	1.86	0.57
1:X:11:PRO:HB3	1:X:43:GLN:HB3	1.90	0.57
1:I:151:LEU:HD11	1:J:62:LEU:HD13	2.04	0.57
1:1:143:THR:HG22	1:2:55:PRO:HB2	1.84	0.57
1:U:55:PRO:HB2	1:Y:143:THR:HG22	1.90	0.57
1:G:143:THR:HG22	1:H:55:PRO:HB2	2.07	0.57
1:L:151:LEU:HD11	1:M:62:LEU:HD13	1.86	0.57
1:2:64:THR:HB	1:2:76:VAL:HG11	1.86	0.57
1:G:24:GLN:HA	1:G:27:ASN:HB2	1.87	0.57
1:1:37:ALA:HB1	1:1:142:LEU:HD13	1.87	0.57
1:I:122:ILE:H	1:I:122:ILE:CD1	2.11	0.57
1:E:122:ILE:H	1:E:122:ILE:CD1	2.16	0.57
1:4:122:ILE:H	1:4:122:ILE:CD1	2.16	0.57
1:X:77:VAL:HG12	1:X:79:LEU:HD13	1.86	0.57
1:F:153:ALA:O	1:F:154:ILE:HG12	2.04	0.57
1:P:24:GLN:HA	1:P:27:ASN:HB2	1.86	0.57
1:X:127:GLU:HA	1:X:131:THR:OG1	2.05	0.57
1:7:146:GLU:O	1:7:150:VAL:HG23	2.04	0.57
1:M:2:ASN:HD21	1:N:50:THR:HG22	1.69	0.57
1:M:122:ILE:N	1:M:122:ILE:HD12	2.12	0.57
1:7:122:ILE:CD1	1:7:122:ILE:H	2.12	0.57
1:V:24:GLN:HA	1:V:27:ASN:HB2	1.94	0.57
1:X:56:GLY:O	1:X:59:GLU:HG2	2.10	0.57
1:M:56:GLY:O	1:M:59:GLU:HG2	2.20	0.57
1:B:11:PRO:HB3	1:B:43:GLN:HB3	1.90	0.57
1:5:5:LYS:HE2	1:6:21:ARG:NH2	2.16	0.57
1:Q:19:ILE:HG12	1:Q:79:LEU:HB2	1.86	0.57
1:L:11:PRO:HB3	1:L:43:GLN:HB3	1.92	0.57
1:4:37:ALA:HB1	1:4:142:LEU:HD13	1.87	0.57
1:Z:11:PRO:HB3	1:Z:43:GLN:HB3	1.87	0.57
1:U:5:LYS:CE	1:V:21:ARG:HH22	2.30	0.57
1:U:77:VAL:HG12	1:U:79:LEU:HD13	1.87	0.57
1:I:41:ILE:HA	1:O:7:ASN:ND2	2.20	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:S:37:ALA:HB1	1:S:142:LEU:HD13	1.86	0.57
1:O:64:THR:HB	1:O:76:VAL:HG11	1.88	0.57
1:R:41:ILE:HA	1:W:7:ASN:ND2	99.22	0.57
1:T:127:GLU:HA	1:T:131:THR:OG1	2.06	0.57
1:C:127:GLU:HA	1:C:131:THR:OG1	2.04	0.57
1:Q:42:GLY:HA3	1:Q:145:LEU:HD13	1.86	0.57
1:G:77:VAL:HG12	1:G:79:LEU:HD13	1.86	0.57
1:F:50:THR:HG22	1:J:2:ASN:HD21	1.68	0.57
1:B:2:ASN:H	1:B:2:ASN:HD22	1.57	0.57
1:9:146:GLU:O	1:9:150:VAL:HG23	2.05	0.57
1:O:127:GLU:HA	1:O:131:THR:OG1	2.04	0.57
1:I:114:PHE:HB2	1:J:58:TYR:CE2	2.40	0.56
1:M:11:PRO:HB3	1:M:43:GLN:HB3	1.89	0.56
1:W:16:ALA:HB1	1:W:67:LEU:HD13	1.86	0.56
1:O:56:GLY:O	1:O:59:GLU:HG2	2.13	0.56
1:N:127:GLU:HA	1:N:131:THR:OG1	2.05	0.56
1:U:41:ILE:HA	1:V:7:ASN:ND2	48.03	0.56
1:Y:37:ALA:HB1	1:Y:142:LEU:HD13	1.87	0.56
1:I:56:GLY:O	1:I:59:GLU:HG2	2.05	0.56
1:6:122:ILE:HD12	1:6:122:ILE:N	2.11	0.56
1:E:24:GLN:HA	1:E:27:ASN:HB2	1.87	0.56
1:R:37:ALA:HB1	1:R:142:LEU:HD13	1.85	0.56
1:E:127:GLU:HA	1:E:131:THR:OG1	2.05	0.56
1:Q:119:THR:HG22	1:R:87:THR:HG21	1.87	0.56
1:A:41:ILE:HA	1:L:7:ASN:ND2	74.40	0.56
1:R:153:ALA:O	1:R:154:ILE:HG12	2.18	0.56
1:K:46:ASP:O	1:O:1:MET:HA	2.12	0.56
1:N:42:GLY:HA3	1:N:145:LEU:HD13	1.95	0.56
1:E:11:PRO:HB3	1:E:43:GLN:HB3	1.88	0.56
1:K:67:LEU:HD21	1:O:154:ILE:HD13	1.90	0.56
1:2:127:GLU:HA	1:2:131:THR:OG1	2.05	0.56
1:S:64:THR:HB	1:S:76:VAL:HG11	1.88	0.56
1:R:5:LYS:CE	1:S:21:ARG:HH22	2.17	0.56
1:Y:23:ASN:HA	1:4:131:THR:HA	1.86	0.56
1:Q:77:VAL:HG12	1:Q:79:LEU:HD13	1.88	0.56
1:G:118:THR:O	1:H:87:THR:HB	2.42	0.56
1:T:77:VAL:HG12	1:T:79:LEU:HD13	1.87	0.56
1:3:77:VAL:HG12	1:3:79:LEU:HD13	1.87	0.56
1:M:143:THR:HG22	1:N:55:PRO:HB2	1.87	0.56
1:D:122:ILE:HD12	1:D:122:ILE:N	2.13	0.56
1:G:122:ILE:H	1:G:122:ILE:CD1	2.13	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:42:GLY:HA3	1:C:145:LEU:HD13	1.91	0.56
1:L:127:GLU:HA	1:L:131:THR:OG1	2.06	0.56
1:U:151:LEU:HD11	1:V:62:LEU:HD13	1.97	0.56
1:R:19:ILE:HG21	1:R:53:TRP:CZ3	2.41	0.56
1:Y:16:ALA:HB1	1:Y:67:LEU:HD13	1.88	0.56
1:W:37:ALA:HB1	1:W:142:LEU:HD13	1.89	0.56
1:6:64:THR:HB	1:6:76:VAL:HG11	1.88	0.56
1:R:122:ILE:HD12	1:R:122:ILE:N	2.19	0.56
1:Y:127:GLU:HA	1:Y:131:THR:OG1	2.05	0.56
1:F:122:ILE:H	1:F:122:ILE:CD1	2.17	0.56
1:S:143:THR:HG22	1:T:55:PRO:HB2	2.03	0.56
1:B:7:ASN:ND2	1:K:41:ILE:HA	118.69	0.56
1:I:77:VAL:HG12	1:I:79:LEU:HD13	1.87	0.56
1:E:56:GLY:O	1:E:59:GLU:HG2	2.10	0.56
1:B:85:GLY:HA3	2:B:157:SO4:O1	2.76	0.56
1:F:16:ALA:HB1	1:F:67:LEU:HD13	1.91	0.56
1:1:77:VAL:HG12	1:1:79:LEU:HD13	1.87	0.56
1:U:64:THR:HB	1:U:76:VAL:HG11	1.88	0.56
1:R:131:THR:HA	1:X:23:ASN:OD1	75.28	0.56
1:C:5:LYS:HE2	1:D:21:ARG:NH2	2.14	0.56
1:L:23:ASN:HA	1:L:131:THR:HA	17.50	0.56
1:T:11:PRO:HB3	1:T:43:GLN:HB3	1.87	0.56
1:K:2:ASN:HD21	1:L:50:THR:HG22	1.72	0.56
1:7:37:ALA:HB1	1:7:142:LEU:HD13	1.87	0.56
1:P:153:ALA:O	1:P:154:ILE:HG12	2.18	0.56
1:6:127:GLU:HA	1:6:131:THR:OG1	2.06	0.56
1:D:56:GLY:O	1:D:59:GLU:HG2	2.18	0.56
1:H:11:PRO:HB3	1:H:43:GLN:HB3	1.87	0.56
1:9:37:ALA:HB1	1:9:142:LEU:HD13	1.87	0.56
1:A:64:THR:HB	1:A:76:VAL:HG11	1.88	0.56
1:6:24:GLN:HA	1:6:27:ASN:HB2	1.88	0.56
1:P:77:VAL:HG12	1:P:79:LEU:HD13	1.88	0.56
1:V:16:ALA:HB1	1:V:67:LEU:HD13	1.90	0.56
1:I:54:VAL:HB	1:I:55:PRO:HD2	1.95	0.56
1:Q:56:GLY:O	1:Q:59:GLU:HG2	2.06	0.56
1:5:37:ALA:HB1	1:5:142:LEU:HD13	1.86	0.56
1:L:85:GLY:HA3	2:L:157:SO4:O2	2.05	0.56
1:R:122:ILE:HD13	1:Z:123:GLU:OE2	70.04	0.55
1:N:122:ILE:HD13	1:W:123:GLU:OE2	53.46	0.55
1:P:122:ILE:HD12	1:P:122:ILE:N	2.20	0.55
1:A:21:ARG:NH2	1:E:5:LYS:HE2	2.19	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:2:ASN:HD21	1:H:50:THR:CG2	2.18	0.55
1:R:22:PHE:HE1	1:Z:131:THR:O	87.78	0.55
1:L:153:ALA:O	1:L:154:ILE:HG12	2.17	0.55
1:V:77:VAL:HG12	1:V:79:LEU:HD13	1.93	0.55
1:O:42:GLY:HA3	1:O:145:LEU:HD13	1.93	0.55
1:8:37:ALA:HB1	1:8:142:LEU:HD13	1.86	0.55
1:M:16:ALA:HB1	1:M:67:LEU:HD13	1.91	0.55
1:X:153:ALA:O	1:X:154:ILE:HG12	2.05	0.55
1:M:119:THR:HG22	1:N:87:THR:HG21	2.15	0.55
1:B:119:THR:HG22	1:C:87:THR:HG21	1.88	0.55
1:M:126:ILE:HG23	1:X:25:PHE:CE1	81.46	0.55
1:F:64:THR:HB	1:F:76:VAL:HG11	1.91	0.55
1:K:19:ILE:HG12	1:K:79:LEU:HB2	1.95	0.55
1:W:11:PRO:HB3	1:W:43:GLN:HB3	1.90	0.55
1:6:77:VAL:HG12	1:6:79:LEU:HD13	1.88	0.55
1:C:11:PRO:HB3	1:C:43:GLN:HB3	1.97	0.55
1:2:16:ALA:HB1	1:2:67:LEU:HD13	1.87	0.55
1:F:77:VAL:HG12	1:F:79:LEU:HD13	1.88	0.55
1:O:23:ASN:HA	1:T:131:THR:HA	90.52	0.55
1:N:143:THR:HG22	1:O:55:PRO:HB2	1.97	0.55
1:I:16:ALA:HB1	1:I:67:LEU:HD13	1.88	0.55
1:R:2:ASN:HD21	1:S:50:THR:HG22	1.72	0.55
1:S:42:GLY:HA3	1:S:145:LEU:HD13	1.91	0.55
1:S:56:GLY:O	1:S:59:GLU:HG2	2.06	0.55
1:G:122:ILE:N	1:G:122:ILE:HD12	2.18	0.55
1:V:127:GLU:HA	1:V:131:THR:OG1	2.09	0.55
1:B:127:GLU:HA	1:B:131:THR:OG1	2.08	0.55
1:Q:11:PRO:HB3	1:Q:43:GLN:HB3	1.88	0.55
1:I:127:GLU:HA	1:I:131:THR:OG1	2.06	0.55
1:Z:2:ASN:HD21	1:1:50:THR:HG22	1.72	0.55
1:T:24:GLN:HA	1:T:27:ASN:HB2	1.88	0.55
1:Z:24:GLN:HA	1:Z:27:ASN:HB2	1.88	0.55
1:L:24:GLN:HA	1:L:27:ASN:HB2	1.88	0.55
1:K:16:ALA:HB1	1:K:67:LEU:HD13	1.88	0.55
1:J:11:PRO:HB3	1:J:43:GLN:HB3	1.89	0.55
1:V:42:GLY:HA3	1:V:145:LEU:HD13	1.89	0.55
1:Y:122:ILE:H	1:Y:122:ILE:CD1	2.13	0.55
1:1:24:GLN:HA	1:1:27:ASN:HB2	1.89	0.55
1:3:119:THR:HG22	1:4:87:THR:HG21	1.88	0.55
1:O:54:VAL:HB	1:O:55:PRO:HD2	1.91	0.55
1:Q:16:ALA:HB1	1:Q:67:LEU:HD13	1.88	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:V:19:ILE:HG12	1:V:79:LEU:HB2	1.91	0.55
1:P:16:ALA:HB1	1:P:67:LEU:HD13	1.89	0.55
1:O:85:GLY:HA3	2:O:157:SO4:O1	2.07	0.55
1:W:119:THR:HG22	1:X:87:THR:HG21	1.89	0.55
1:R:16:ALA:HB1	1:R:67:LEU:HD13	1.89	0.55
1:V:64:THR:HB	1:V:76:VAL:HG11	1.89	0.55
1:P:11:PRO:HB3	1:P:43:GLN:HB3	1.88	0.55
1:A:58:TYR:CE2	1:E:114:PHE:HB2	2.46	0.55
1:I:42:GLY:HA3	1:I:145:LEU:HD13	1.91	0.55
1:7:127:GLU:HA	1:7:131:THR:OG1	2.07	0.55
1:G:154:ILE:HD13	1:H:67:LEU:CD2	2.62	0.55
1:P:127:GLU:HA	1:P:131:THR:OG1	2.07	0.55
2:3:157:SO4:O4	1:4:85:GLY:HA3	2.07	0.55
1:3:16:ALA:HB1	1:3:67:LEU:HD13	1.89	0.55
1:F:24:GLN:HA	1:F:27:ASN:HB2	1.88	0.54
1:G:5:LYS:HE2	1:H:21:ARG:NH2	2.18	0.54
1:M:19:ILE:HG21	1:M:53:TRP:CZ3	2.51	0.54
1:N:19:ILE:HG12	1:N:79:LEU:HB2	1.90	0.54
1:K:21:ARG:O	1:K:24:GLN:HB3	2.07	0.54
1:W:54:VAL:HB	1:W:55:PRO:HD2	1.89	0.54
1:I:39:THR:OG1	1:I:40:ARG:N	2.39	0.54
1:X:42:GLY:HA3	1:X:145:LEU:HD13	1.89	0.54
1:Q:143:THR:HG22	1:R:55:PRO:HB2	1.88	0.54
1:G:55:PRO:HG2	1:G:59:GLU:HG3	1.89	0.54
1:9:77:VAL:HG12	1:9:79:LEU:HD13	1.89	0.54
1:V:122:ILE:N	1:V:122:ILE:HD12	2.15	0.54
1:M:19:ILE:HG12	1:M:79:LEU:HB2	1.88	0.54
1:U:53:TRP:CD1	1:Y:5:LYS:HE3	2.42	0.54
1:V:151:LEU:HD11	1:W:62:LEU:HD13	1.95	0.54
1:8:11:PRO:HB3	1:8:43:GLN:HB3	1.89	0.54
1:J:56:GLY:O	1:J:59:GLU:HG2	2.08	0.54
1:G:127:GLU:HA	1:G:131:THR:OG1	2.07	0.54
1:S:24:GLN:HA	1:S:27:ASN:HB2	1.91	0.54
1:Y:42:GLY:HA3	1:Y:145:LEU:HD13	1.89	0.54
1:M:55:PRO:HG2	1:M:59:GLU:HG3	1.90	0.54
1:R:42:GLY:HA3	1:R:145:LEU:HD13	1.91	0.54
1:Z:114:PHE:HB2	1:1:58:TYR:CE2	2.42	0.54
1:F:42:GLY:HA3	1:F:145:LEU:HD13	1.92	0.54
1:9:2:ASN:HD22	1:9:2:ASN:H	1.54	0.54
1:D:19:ILE:HG12	1:D:79:LEU:HB2	1.91	0.54
1:C:24:GLN:HA	1:C:27:ASN:HB2	1.90	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:U:42:GLY:HA3	1:U:145:LEU:HD13	1.88	0.54
1:D:54:VAL:HB	1:D:55:PRO:HD2	1.95	0.54
1:F:55:PRO:HG2	1:F:59:GLU:HG3	1.89	0.54
1:8:16:ALA:HB1	1:8:67:LEU:HD13	1.89	0.54
1:J:77:VAL:HG12	1:J:79:LEU:HD13	1.89	0.54
1:M:127:GLU:HA	1:M:131:THR:OG1	2.07	0.54
1:A:7:ASN:ND2	1:V:41:ILE:HA	37.28	0.54
1:K:55:PRO:HB2	1:O:143:THR:HG22	1.91	0.54
1:U:56:GLY:O	1:U:59:GLU:HG2	2.08	0.54
1:8:2:ASN:H	1:8:2:ASN:HD22	1.55	0.54
1:A:16:ALA:HB1	1:A:67:LEU:HD13	1.94	0.54
1:Z:19:ILE:HG12	1:Z:79:LEU:HB2	1.89	0.54
1:S:77:VAL:HG12	1:S:79:LEU:HD13	1.93	0.54
1:N:111:PRO:HG3	1:N:151:LEU:HD12	2.02	0.54
1:K:77:VAL:HG12	1:K:79:LEU:HD13	1.93	0.54
1:2:23:ASN:HA	1:5:131:THR:HA	1.88	0.54
1:Y:24:GLN:HA	1:Y:27:ASN:HB2	1.90	0.54
1:K:127:GLU:HA	1:K:131:THR:OG1	2.09	0.54
1:Q:19:ILE:HG21	1:Q:53:TRP:CZ3	2.42	0.54
1:7:56:GLY:O	1:7:59:GLU:HG2	2.08	0.54
1:C:16:ALA:HB1	1:C:67:LEU:HD13	1.93	0.54
1:C:19:ILE:HG21	1:C:53:TRP:CZ3	2.50	0.54
1:V:55:PRO:HG2	1:V:59:GLU:HG3	1.93	0.54
1:7:42:GLY:HA3	1:7:145:LEU:HD13	1.90	0.54
1:Z:37:ALA:HB1	1:Z:142:LEU:HD13	1.89	0.54
1:J:16:ALA:HB1	1:J:67:LEU:HD13	1.91	0.54
1:8:127:GLU:HA	1:8:131:THR:OG1	2.07	0.54
1:9:21:ARG:O	1:9:24:GLN:HB3	2.08	0.54
1:Y:77:VAL:HG12	1:Y:79:LEU:HD13	1.89	0.54
1:N:131:THR:HA	1:P:23:ASN:HA	90.00	0.54
1:L:42:GLY:HA3	1:L:145:LEU:HD13	1.90	0.54
1:C:77:VAL:HG12	1:C:79:LEU:HD13	1.89	0.54
1:H:42:GLY:HA3	1:H:145:LEU:HD13	1.92	0.54
1:X:143:THR:HG22	1:Y:55:PRO:HB2	1.94	0.54
1:Z:122:ILE:CD1	1:Z:122:ILE:H	2.08	0.54
1:Y:19:ILE:HG12	1:Y:79:LEU:HB2	1.90	0.54
1:2:2:ASN:HD21	1:3:50:THR:HG22	1.73	0.54
1:T:40:ARG:O	1:3:7:ASN:ND2	2.41	0.54
1:S:119:THR:HG22	1:T:87:THR:HG21	2.08	0.54
1:4:77:VAL:HG12	1:4:79:LEU:HD13	1.90	0.54
1:Q:2:ASN:HD22	1:Q:2:ASN:H	1.58	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:24:GLN:HA	1:H:27:ASN:HB2	1.89	0.53
1:Q:7:ASN:HD22	1:Z:41:ILE:HD13	117.79	0.53
1:P:87:THR:HB	1:T:118:THR:O	2.07	0.53
1:U:154:ILE:HD13	1:V:67:LEU:CD2	2.63	0.53
1:W:143:THR:HG22	1:X:55:PRO:HB2	1.90	0.53
1:L:16:ALA:HB1	1:L:67:LEU:HD13	1.90	0.53
1:I:21:ARG:O	1:I:24:GLN:HB3	2.08	0.53
1:W:127:GLU:HA	1:W:131:THR:OG1	2.08	0.53
1:R:2:ASN:HD22	1:R:2:ASN:H	1.56	0.53
1:N:16:ALA:HB1	1:N:67:LEU:HD13	1.90	0.53
1:A:11:PRO:HB3	1:A:43:GLN:HB3	1.90	0.53
1:O:16:ALA:HB1	1:O:67:LEU:HD13	1.90	0.53
1:8:122:ILE:H	1:8:122:ILE:CD1	2.13	0.53
1:M:77:VAL:HG12	1:M:79:LEU:HD13	1.90	0.53
1:U:19:ILE:HG12	1:U:79:LEU:HB2	1.90	0.53
1:8:114:PHE:HB2	1:9:58:TYR:CE2	2.43	0.53
1:N:11:PRO:HB3	1:N:43:GLN:HB3	1.90	0.53
1:X:1:MET:HB2	1:Y:46:ASP:O	2.24	0.53
1:B:55:PRO:HG2	1:B:59:GLU:HG3	1.91	0.53
1:X:16:ALA:HB1	1:X:67:LEU:HD13	1.94	0.53
1:B:114:PHE:HB2	1:C:58:TYR:CE2	2.46	0.53
1:V:5:LYS:CE	1:W:21:ARG:HH22	2.13	0.53
1:U:19:ILE:HG21	1:U:53:TRP:CZ3	2.43	0.53
1:V:2:ASN:HD22	1:V:2:ASN:H	1.61	0.53
1:Z:127:GLU:HA	1:Z:131:THR:OG1	2.09	0.53
1:T:42:GLY:HA3	1:T:145:LEU:HD13	1.91	0.53
1:I:19:ILE:HG12	1:I:79:LEU:HB2	1.90	0.53
1:8:24:GLN:HA	1:8:27:ASN:HB2	1.90	0.53
1:Q:141:ALA:O	1:Q:144:ALA:HB3	2.08	0.53
1:Z:55:PRO:HB2	1:4:143:THR:HG22	1.91	0.53
1:5:16:ALA:HB1	1:5:67:LEU:HD13	1.91	0.53
1:R:130:GLY:O	1:X:23:ASN:HA	79.41	0.53
1:W:24:GLN:HA	1:W:27:ASN:HB2	1.90	0.53
1:R:5:LYS:HE2	1:S:21:ARG:NH2	2.18	0.53
1:X:122:ILE:HD12	1:X:122:ILE:N	2.18	0.53
1:B:77:VAL:HG12	1:B:79:LEU:HD13	1.92	0.53
1:3:127:GLU:HA	1:3:131:THR:OG1	2.08	0.53
1:7:16:ALA:HB1	1:7:67:LEU:HD13	1.90	0.53
1:D:2:ASN:HD22	1:D:2:ASN:H	1.57	0.53
1:9:55:PRO:HG2	1:9:59:GLU:HG3	1.90	0.53
1:V:39:THR:OG1	1:V:40:ARG:N	2.41	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:77:VAL:HG12	1:D:79:LEU:HD13	1.92	0.53
1:9:122:ILE:H	1:9:122:ILE:CD1	2.11	0.53
1:8:5:LYS:HE2	1:9:21:ARG:NH2	2.21	0.53
1:B:19:ILE:HG21	1:B:53:TRP:CZ3	2.51	0.53
1:5:21:ARG:NH2	1:9:5:LYS:HE2	2.19	0.53
1:T:41:ILE:HA	1:3:7:ASN:ND2	2.23	0.53
1:N:56:GLY:O	1:N:59:GLU:HG2	2.09	0.53
1:G:54:VAL:HB	1:G:55:PRO:HD2	1.93	0.53
1:O:39:THR:OG1	1:O:40:ARG:N	2.51	0.53
1:L:2:ASN:H	1:L:2:ASN:HD22	1.62	0.53
1:P:19:ILE:HG21	1:P:53:TRP:CZ3	2.49	0.53
1:L:154:ILE:HD13	1:M:67:LEU:HD21	1.98	0.53
1:P:42:GLY:HA3	1:P:145:LEU:HD13	1.89	0.53
1:M:42:GLY:HA3	1:M:145:LEU:HD13	1.94	0.53
1:G:11:PRO:HB3	1:G:43:GLN:HB3	1.90	0.53
1:2:114:PHE:HB2	1:3:58:TYR:CE2	2.44	0.53
1:1:2:ASN:HD22	1:1:2:ASN:H	1.56	0.53
1:D:64:THR:HB	1:D:76:VAL:HG11	1.89	0.53
1:D:21:ARG:O	1:D:24:GLN:HB3	2.11	0.53
1:9:127:GLU:HA	1:9:131:THR:OG1	2.08	0.53
1:R:151:LEU:HD11	1:S:62:LEU:HD13	1.90	0.53
1:B:56:GLY:O	1:B:59:GLU:HG2	2.20	0.53
1:L:64:THR:HB	1:L:76:VAL:HG11	1.90	0.53
1:F:19:ILE:HG12	1:F:79:LEU:HB2	1.93	0.53
1:H:77:VAL:HG12	1:H:79:LEU:HD13	1.91	0.53
1:P:55:PRO:HG2	1:P:59:GLU:HG3	1.91	0.53
1:W:56:GLY:O	1:W:59:GLU:HG2	2.08	0.53
1:4:19:ILE:HG21	1:4:53:TRP:CZ3	2.44	0.53
1:M:151:LEU:O	1:M:154:ILE:HG13	2.21	0.53
1:5:77:VAL:HG12	1:5:79:LEU:HD13	1.91	0.53
1:4:16:ALA:HB1	1:4:67:LEU:HD13	1.89	0.53
1:U:122:ILE:CD1	1:U:122:ILE:H	2.14	0.53
1:O:153:ALA:O	1:O:154:ILE:HG12	2.09	0.53
1:Z:54:VAL:HB	1:Z:55:PRO:HD2	1.90	0.53
1:Z:55:PRO:HG2	1:Z:59:GLU:HG3	1.91	0.53
1:A:19:ILE:HG21	1:A:53:TRP:CZ3	2.49	0.53
1:D:153:ALA:O	1:D:154:ILE:HG12	2.12	0.52
1:T:54:VAL:HB	1:T:55:PRO:HD2	1.97	0.52
1:F:21:ARG:HH22	1:J:5:LYS:CE	2.20	0.52
1:N:24:GLN:HA	1:N:27:ASN:HB2	1.91	0.52
1:K:24:GLN:HA	1:K:27:ASN:HB2	1.92	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:Y:21:ARG:O	1:Y:24:GLN:HB3	2.14	0.52
1:E:19:ILE:HG12	1:E:79:LEU:HB2	1.94	0.52
1:B:42:GLY:HA3	1:B:145:LEU:HD13	1.90	0.52
1:E:41:ILE:HA	1:N:7:ASN:ND2	2.24	0.52
1:6:2:ASN:HD22	1:6:2:ASN:H	1.57	0.52
1:Z:42:GLY:HA3	1:Z:145:LEU:HD13	1.91	0.52
1:K:39:THR:OG1	1:K:40:ARG:N	2.42	0.52
1:T:111:PRO:HG3	1:T:151:LEU:HD12	1.92	0.52
1:B:31:LEU:O	1:B:35:VAL:HG23	2.22	0.52
1:M:31:LEU:O	1:M:35:VAL:HG23	2.09	0.52
1:G:21:ARG:O	1:G:24:GLN:HB3	2.12	0.52
1:I:2:ASN:HD22	1:I:2:ASN:H	1.62	0.52
1:D:16:ALA:HB1	1:D:67:LEU:HD13	1.91	0.52
1:E:111:PRO:HG3	1:E:151:LEU:HD12	2.00	0.52
1:B:122:ILE:CD1	1:B:122:ILE:H	2.15	0.52
1:C:5:LYS:CE	1:D:21:ARG:HH22	2.16	0.52
1:D:24:GLN:HA	1:D:27:ASN:HB2	1.94	0.52
1:U:50:THR:CG2	1:Y:2:ASN:HD21	2.21	0.52
1:U:21:ARG:O	1:U:24:GLN:HB3	2.10	0.52
1:W:55:PRO:HG2	1:W:59:GLU:HG3	1.91	0.52
1:V:56:GLY:O	1:V:59:GLU:HG2	2.16	0.52
1:Q:54:VAL:HB	1:Q:55:PRO:HD2	1.91	0.52
1:J:19:ILE:HG12	1:J:79:LEU:HB2	1.91	0.52
1:L:119:THR:HG22	1:M:87:THR:HG21	1.91	0.52
1:7:77:VAL:HG12	1:7:79:LEU:HD13	1.90	0.52
1:Q:122:ILE:HD12	1:Q:122:ILE:N	2.19	0.52
1:I:24:GLN:HA	1:I:27:ASN:HB2	1.91	0.52
1:K:21:ARG:CZ	1:X:40:ARG:CZ	126.56	0.52
1:G:7:ASN:ND2	1:P:41:ILE:HA	50.07	0.52
1:H:55:PRO:HG2	1:H:59:GLU:HG3	1.98	0.52
1:M:2:ASN:HD22	1:M:2:ASN:H	1.58	0.52
1:D:7:ASN:ND2	1:J:41:ILE:HA	2.24	0.52
1:7:21:ARG:O	1:7:24:GLN:HB3	2.10	0.52
1:O:21:ARG:O	1:O:24:GLN:HB3	2.10	0.52
1:O:24:GLN:HA	1:O:27:ASN:HB2	1.98	0.52
1:U:119:THR:HA	1:V:87:THR:HB	1.91	0.52
1:A:7:ASN:ND2	1:L:41:ILE:HA	87.63	0.52
1:Z:77:VAL:HG12	1:Z:79:LEU:HD13	1.91	0.52
1:1:2:ASN:HD21	1:2:50:THR:HG22	1.73	0.52
1:U:24:GLN:HA	1:U:27:ASN:HB2	1.91	0.52
1:H:56:GLY:O	1:H:59:GLU:HG2	2.17	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:V:118:THR:O	1:W:87:THR:HB	2.30	0.52
1:4:42:GLY:HA3	1:4:145:LEU:HD13	1.92	0.52
1:B:122:ILE:HD12	1:B:122:ILE:N	2.18	0.52
1:D:5:LYS:HE2	1:E:21:ARG:NH2	2.19	0.52
1:X:39:THR:OG1	1:X:40:ARG:N	2.51	0.52
1:U:153:ALA:O	1:U:154:ILE:HG12	2.19	0.52
1:3:19:ILE:HG12	1:3:79:LEU:HB2	1.92	0.52
1:S:122:ILE:HD12	1:S:122:ILE:N	2.19	0.52
1:V:119:THR:HG22	1:W:87:THR:HG21	2.14	0.52
1:A:56:GLY:O	1:A:59:GLU:HG2	2.12	0.52
1:9:64:THR:HB	1:9:76:VAL:HG11	1.91	0.52
1:K:122:ILE:HD12	1:K:122:ILE:N	2.19	0.52
1:A:7:ASN:ND2	1:L:40:ARG:O	89.97	0.52
1:W:19:ILE:HG12	1:W:79:LEU:HB2	1.92	0.52
1:6:143:THR:HG22	1:7:55:PRO:HB2	1.92	0.52
1:7:19:ILE:HG21	1:7:53:TRP:CZ3	2.44	0.52
1:8:39:THR:OG1	1:8:40:ARG:N	2.42	0.52
1:T:55:PRO:HG2	1:T:59:GLU:HG3	1.92	0.52
1:C:122:ILE:CD1	1:C:122:ILE:H	2.15	0.52
1:B:5:LYS:HE2	1:C:21:ARG:NH2	2.19	0.52
1:K:21:ARG:NH2	1:O:5:LYS:HE2	2.25	0.52
1:G:2:ASN:HD22	1:G:2:ASN:H	1.63	0.52
1:L:56:GLY:O	1:L:59:GLU:HG2	2.19	0.52
1:3:39:THR:OG1	1:3:40:ARG:N	2.43	0.52
1:X:2:ASN:H	1:X:2:ASN:HD22	1.63	0.52
1:3:114:PHE:HB2	1:4:58:TYR:CE2	2.45	0.52
1:3:42:GLY:HA3	1:3:145:LEU:HD13	1.92	0.52
1:I:122:ILE:HD12	1:I:122:ILE:N	2.17	0.51
1:F:21:ARG:O	1:F:24:GLN:HB3	2.14	0.51
1:2:5:LYS:HE2	1:3:21:ARG:NH2	2.21	0.51
1:N:19:ILE:HG12	1:N:79:LEU:CB	2.40	0.51
1:S:2:ASN:H	1:S:2:ASN:HD22	1.58	0.51
1:7:2:ASN:HD22	1:7:2:ASN:H	1.57	0.51
1:8:122:ILE:N	1:8:122:ILE:HD12	2.18	0.51
1:M:54:VAL:HB	1:M:55:PRO:HD2	1.92	0.51
1:D:55:PRO:HG2	1:D:59:GLU:HG3	1.92	0.51
1:V:31:LEU:O	1:V:34:ALA:N	2.44	0.51
1:C:111:PRO:HG3	1:C:151:LEU:HD12	1.91	0.51
1:L:77:VAL:HG12	1:L:79:LEU:HD13	1.92	0.51
1:M:24:GLN:HA	1:M:27:ASN:HB2	1.92	0.51
1:Q:21:ARG:O	1:Q:24:GLN:HB3	2.14	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:K:19:ILE:HG21	1:K:53:TRP:CZ3	2.45	0.51
1:N:2:ASN:HD22	1:N:2:ASN:H	1.59	0.51
1:E:2:ASN:H	1:E:2:ASN:HD22	1.59	0.51
1:A:42:GLY:HA3	1:A:145:LEU:HD13	1.94	0.51
1:A:62:LEU:HD13	1:E:151:LEU:HD11	1.92	0.51
1:J:42:GLY:HA3	1:J:145:LEU:HD13	1.93	0.51
1:E:123:GLU:OE2	1:O:122:ILE:HD13	2.11	0.51
1:X:21:ARG:O	1:X:24:GLN:HB3	2.13	0.51
1:G:42:GLY:HA3	1:G:145:LEU:HD13	1.92	0.51
1:I:19:ILE:HG12	1:I:79:LEU:HB2	1.91	0.51
1:6:19:ILE:HG21	1:6:53:TRP:CZ3	2.45	0.51
1:S:19:ILE:HG12	1:S:79:LEU:HB2	1.96	0.51
1:U:16:ALA:HB1	1:U:67:LEU:HD13	1.93	0.51
1:G:64:THR:HB	1:G:76:VAL:HG11	1.92	0.51
1:N:123:GLU:OE2	1:P:122:ILE:HD13	74.67	0.51
1:I:27:ASN:HA	1:I:30:LEU:HB2	1.92	0.51
1:4:21:ARG:O	1:4:24:GLN:HB3	2.11	0.51
1:H:119:THR:HG22	1:I:87:THR:HG21	1.92	0.51
1:J:2:ASN:H	1:J:2:ASN:HD22	1.61	0.51
1:U:62:LEU:HD13	1:Y:151:LEU:HD11	1.92	0.51
1:6:16:ALA:HB1	1:6:67:LEU:HD13	1.92	0.51
1:F:118:THR:O	1:G:87:THR:HB	2.11	0.51
1:J:24:GLN:HA	1:J:27:ASN:HB2	1.93	0.51
1:M:64:THR:HB	1:M:76:VAL:HG11	1.91	0.51
1:N:122:ILE:HD12	1:N:122:ILE:N	2.20	0.51
1:T:56:GLY:O	1:T:59:GLU:HG2	2.10	0.51
1:W:64:THR:HB	1:W:76:VAL:HG11	1.94	0.51
1:B:24:GLN:HA	1:B:27:ASN:HB2	1.96	0.51
1:W:5:LYS:HE2	1:X:21:ARG:NH2	2.21	0.51
1:R:111:PRO:HG3	1:R:151:LEU:HD12	1.92	0.51
1:A:54:VAL:HB	1:A:55:PRO:HD2	1.97	0.51
1:6:42:GLY:HA3	1:6:145:LEU:HD13	1.93	0.51
1:K:42:GLY:HA3	1:K:145:LEU:HD13	1.92	0.51
1:Z:16:ALA:HB1	1:Z:67:LEU:HD13	1.93	0.51
1:E:16:ALA:HB1	1:E:67:LEU:HD13	1.93	0.51
1:O:122:ILE:HD12	1:O:122:ILE:N	2.20	0.51
1:S:153:ALA:O	1:S:154:ILE:HG12	2.11	0.51
1:X:122:ILE:CD1	1:X:122:ILE:H	2.14	0.51
1:A:21:ARG:O	1:A:24:GLN:HB3	2.14	0.51
1:C:21:ARG:O	1:C:24:GLN:HB3	2.12	0.51
1:Z:21:ARG:O	1:Z:24:GLN:HB3	2.11	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:I:55:PRO:HG2	1:I:59:GLU:HG3	1.92	0.51
1:D:42:GLY:HA3	1:D:145:LEU:HD13	1.92	0.51
1:R:118:THR:O	1:S:87:THR:HB	2.12	0.51
1:W:122:ILE:N	1:W:122:ILE:HD12	2.18	0.51
1:A:21:ARG:HH22	1:E:5:LYS:CE	2.25	0.51
1:H:19:ILE:HG21	1:H:53:TRP:CZ3	2.53	0.51
1:X:55:PRO:HG2	1:X:59:GLU:HG3	1.96	0.51
1:2:39:THR:OG1	1:2:40:ARG:N	2.44	0.51
1:A:19:ILE:HG12	1:A:79:LEU:HB2	1.99	0.51
1:2:77:VAL:HG12	1:2:79:LEU:HD13	1.92	0.51
1:D:19:ILE:HG12	1:D:79:LEU:CB	2.48	0.51
1:D:5:LYS:CE	1:E:21:ARG:HH22	2.20	0.51
1:X:19:ILE:HG21	1:X:53:TRP:CZ3	2.46	0.51
1:C:131:THR:HA	1:I:23:ASN:HA	77.22	0.51
1:P:2:ASN:H	1:P:2:ASN:HD22	1.58	0.51
1:G:19:ILE:HG12	1:G:79:LEU:HB2	1.93	0.51
1:B:54:VAL:HB	1:B:55:PRO:HD2	1.93	0.51
1:Y:64:THR:HB	1:Y:76:VAL:HG11	1.93	0.51
1:B:19:ILE:HG12	1:B:79:LEU:HB2	1.92	0.51
1:P:5:LYS:HE2	1:Q:21:ARG:NH2	2.22	0.51
1:X:24:GLN:HA	1:X:27:ASN:HB2	1.94	0.51
1:T:7:ASN:ND2	1:3:41:ILE:HA	2.26	0.51
1:H:54:VAL:HB	1:H:55:PRO:HD2	1.93	0.51
1:U:41:ILE:HD12	1:V:8:VAL:HG22	44.19	0.51
1:Y:55:PRO:HG2	1:Y:59:GLU:HG3	1.93	0.51
1:4:11:PRO:HB3	1:4:43:GLN:HB3	1.93	0.51
1:K:85:GLY:HA3	2:K:157:SO4:O4	4.45	0.51
1:U:1:MET:HA	1:V:46:ASP:O	2.39	0.51
1:5:114:PHE:HB2	1:6:58:TYR:CE2	2.46	0.51
1:5:24:GLN:HA	1:5:27:ASN:HB2	1.92	0.50
1:N:19:ILE:HG21	1:N:53:TRP:CZ3	2.46	0.50
1:T:22:PHE:HD1	1:T:23:ASN:N	2.18	0.50
1:4:2:ASN:H	1:4:2:ASN:HD22	1.58	0.50
1:C:19:ILE:HG12	1:C:79:LEU:HB2	1.91	0.50
1:G:39:THR:OG1	1:G:40:ARG:N	2.47	0.50
1:G:4:ILE:HD12	1:H:52:VAL:HG22	2.26	0.50
1:8:111:PRO:HG3	1:8:151:LEU:HD12	1.93	0.50
1:H:21:ARG:O	1:H:24:GLN:HB3	2.11	0.50
1:T:9:ALA:HA	1:T:42:GLY:HA2	1.93	0.50
1:7:111:PRO:HG3	1:7:151:LEU:HD12	1.92	0.50
1:H:7:ASN:ND2	1:2:41:ILE:HA	99.91	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:5:42:GLY:HA3	1:5:145:LEU:HD13	1.94	0.50
1:1:153:ALA:O	1:1:154:ILE:HG12	2.11	0.50
1:Z:64:THR:HB	1:Z:76:VAL:HG11	1.93	0.50
1:Q:64:THR:HB	1:Q:76:VAL:HG11	1.98	0.50
1:A:24:GLN:HA	1:A:27:ASN:HB2	1.93	0.50
1:F:19:ILE:HG12	1:F:79:LEU:CB	2.44	0.50
1:3:24:GLN:HA	1:3:27:ASN:HB2	1.93	0.50
1:H:19:ILE:HG12	1:H:79:LEU:HB2	1.93	0.50
1:Z:5:LYS:HE2	1:1:21:ARG:NH2	2.21	0.50
1:N:31:LEU:O	1:N:32:ASP:C	2.49	0.50
1:X:19:ILE:HG12	1:X:79:LEU:HB2	1.94	0.50
1:I:19:ILE:HG21	1:I:53:TRP:CZ3	2.49	0.50
1:A:31:LEU:O	1:A:35:VAL:HG23	2.14	0.50
1:9:16:ALA:HB1	1:9:67:LEU:HD13	1.94	0.50
1:B:39:THR:OG1	1:B:40:ARG:N	2.50	0.50
1:V:151:LEU:O	1:V:154:ILE:HG13	2.23	0.50
1:T:19:ILE:HG21	1:T:53:TRP:CZ3	2.47	0.50
1:A:77:VAL:HG12	1:A:79:LEU:HD13	1.93	0.50
1:X:2:ASN:HD21	1:Y:50:THR:HG22	1.76	0.50
1:D:19:ILE:HG21	1:D:53:TRP:CZ3	2.47	0.50
1:G:7:ASN:ND2	1:P:41:ILE:HD13	48.75	0.50
1:B:2:ASN:HD21	1:C:50:THR:HG22	1.77	0.50
1:V:154:ILE:HD13	1:W:67:LEU:HD21	2.11	0.50
1:Q:153:ALA:O	1:Q:154:ILE:HG12	2.12	0.50
1:R:56:GLY:O	1:R:59:GLU:HG2	2.15	0.50
1:8:143:THR:HG22	1:9:55:PRO:HB2	1.93	0.50
1:5:119:THR:HG22	1:6:87:THR:HG21	1.92	0.50
1:9:42:GLY:HA3	1:9:145:LEU:HD13	1.93	0.50
1:1:42:GLY:HA3	1:1:145:LEU:HD13	1.94	0.50
1:3:122:ILE:N	1:3:122:ILE:HD12	2.16	0.50
1:5:5:LYS:CE	1:6:21:ARG:HH22	2.17	0.50
1:U:21:ARG:NH2	1:Y:5:LYS:HE2	2.22	0.50
1:3:2:ASN:H	1:3:2:ASN:HD22	1.60	0.50
1:2:55:PRO:HG2	1:2:59:GLU:HG3	1.93	0.50
1:U:54:VAL:HB	1:U:55:PRO:HD2	1.98	0.50
1:G:119:THR:HG22	1:H:87:THR:HG21	2.17	0.50
1:S:19:ILE:HG21	1:S:53:TRP:CZ3	2.47	0.50
1:V:27:ASN:HA	1:V:30:LEU:HB2	1.94	0.50
1:G:19:ILE:HG21	1:G:53:TRP:CZ3	2.49	0.50
1:M:9:ALA:HA	1:M:42:GLY:HA2	2.04	0.50
1:L:19:ILE:HG12	1:L:79:LEU:HB2	1.95	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:Q:7:ASN:ND2	1:Z:41:ILE:HD13	118.28	0.50
1:L:143:THR:HG22	1:M:55:PRO:HB2	1.96	0.50
1:F:55:PRO:HB2	1:J:143:THR:HG22	1.98	0.50
1:2:2:ASN:H	1:2:2:ASN:HD22	1.60	0.50
1:W:154:ILE:HD13	1:X:67:LEU:HD21	1.93	0.50
1:B:16:ALA:HB1	1:B:67:LEU:HD13	1.94	0.50
1:J:39:THR:OG1	1:J:40:ARG:N	2.44	0.50
1:6:55:PRO:HG2	1:6:59:GLU:HG3	1.93	0.50
1:J:153:ALA:O	1:J:154:ILE:HG12	2.14	0.50
1:H:131:THR:O	1:3:22:PHE:HE1	71.43	0.50
1:V:2:ASN:HD21	1:W:50:THR:CG2	2.24	0.50
1:P:131:THR:HA	1:7:23:ASN:HA	81.90	0.50
1:U:2:ASN:HD22	1:U:2:ASN:H	1.62	0.50
1:2:24:GLN:HA	1:2:27:ASN:HB2	1.93	0.50
1:8:42:GLY:HA3	1:8:145:LEU:HD13	1.94	0.50
1:3:19:ILE:HG21	1:3:53:TRP:CZ3	2.47	0.50
1:F:67:LEU:O	1:F:70:SER:HB3	2.12	0.50
1:S:54:VAL:HB	1:S:55:PRO:HD2	1.94	0.50
1:J:54:VAL:HB	1:J:55:PRO:HD2	1.94	0.50
1:5:143:THR:HG22	1:6:55:PRO:HB2	1.93	0.50
1:T:141:ALA:O	1:T:144:ALA:HB3	2.31	0.50
1:3:55:PRO:HG2	1:3:59:GLU:HG3	1.93	0.50
1:3:56:GLY:O	1:3:59:GLU:HG2	2.12	0.50
1:5:55:PRO:HG2	1:5:59:GLU:HG3	1.92	0.50
1:M:126:ILE:HG23	1:X:25:PHE:CZ	81.63	0.49
1:H:122:ILE:HD12	1:H:122:ILE:N	2.17	0.49
1:R:21:ARG:O	1:R:24:GLN:HB3	2.13	0.49
1:L:21:ARG:O	1:L:24:GLN:HB3	2.15	0.49
1:C:2:ASN:H	1:C:2:ASN:HD22	1.60	0.49
1:H:2:ASN:HD22	1:H:2:ASN:H	1.58	0.49
1:R:19:ILE:HG12	1:R:79:LEU:HB2	1.96	0.49
1:1:19:ILE:HG21	1:1:53:TRP:CZ3	2.46	0.49
1:Q:154:ILE:HD13	1:R:67:LEU:HD21	1.94	0.49
1:9:39:THR:OG1	1:9:40:ARG:N	2.45	0.49
1:A:19:ILE:HD11	1:A:31:LEU:HB2	2.09	0.49
1:Q:39:THR:OG1	1:Q:40:ARG:N	2.45	0.49
1:6:153:ALA:O	1:6:154:ILE:HG12	2.12	0.49
1:R:64:THR:HB	1:R:76:VAL:HG11	1.94	0.49
1:N:21:ARG:O	1:N:24:GLN:HB3	2.15	0.49
1:Q:8:VAL:HG22	1:Z:41:ILE:HD12	115.21	0.49
1:Z:31:LEU:O	1:Z:35:VAL:HG23	2.12	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:K:154:ILE:HD13	1:L:67:LEU:HD21	1.99	0.49
1:B:153:ALA:O	1:B:154:ILE:HG12	2.15	0.49
1:B:143:THR:HG22	1:C:55:PRO:HB2	1.98	0.49
1:E:64:THR:HB	1:E:76:VAL:HG11	1.94	0.49
1:B:18:THR:HG21	1:B:64:THR:HG22	1.94	0.49
1:V:21:ARG:O	1:V:24:GLN:HB3	2.15	0.49
1:3:64:THR:HB	1:3:76:VAL:HG11	1.94	0.49
1:W:19:ILE:HG21	1:W:53:TRP:CZ3	2.47	0.49
1:R:7:ASN:OD1	1:R:9:ALA:HB3	2.12	0.49
1:D:2:ASN:HD21	1:E:50:THR:HG22	1.85	0.49
1:C:54:VAL:HB	1:C:55:PRO:HD2	1.97	0.49
1:8:19:ILE:HG21	1:8:53:TRP:CZ3	2.47	0.49
1:6:39:THR:OG1	1:6:40:ARG:N	2.45	0.49
1:7:39:THR:OG1	1:7:40:ARG:N	2.44	0.49
1:Q:55:PRO:HG2	1:Q:59:GLU:HG3	1.95	0.49
1:V:19:ILE:HG12	1:V:79:LEU:CB	2.43	0.49
1:W:111:PRO:HG3	1:W:151:LEU:HD12	2.04	0.49
1:5:64:THR:HB	1:5:76:VAL:HG11	1.95	0.49
1:4:24:GLN:HA	1:4:27:ASN:HB2	1.93	0.49
1:P:21:ARG:O	1:P:24:GLN:HB3	2.14	0.49
1:U:111:PRO:HG3	1:U:151:LEU:HD12	1.94	0.49
1:T:19:ILE:HG12	1:T:79:LEU:HB2	1.95	0.49
1:8:77:VAL:HG12	1:8:79:LEU:HD13	1.93	0.49
1:U:27:ASN:HA	1:U:30:LEU:HB2	1.95	0.49
1:A:2:ASN:HD21	1:B:50:THR:CG2	2.52	0.49
1:F:2:ASN:H	1:F:2:ASN:HD22	1.60	0.49
1:N:55:PRO:HG2	1:N:59:GLU:HG3	1.94	0.49
1:B:118:THR:O	1:C:87:THR:HB	2.13	0.49
1:4:19:ILE:HG12	1:4:79:LEU:HB2	1.93	0.49
1:9:54:VAL:HB	1:9:55:PRO:HD2	1.95	0.49
1:E:153:ALA:O	1:E:154:ILE:HG12	2.17	0.49
1:8:19:ILE:HG12	1:8:79:LEU:HB2	1.95	0.49
1:K:64:THR:HB	1:K:76:VAL:HG11	1.99	0.49
1:X:31:LEU:O	1:X:32:ASP:C	2.63	0.49
1:E:131:THR:HA	1:9:23:ASN:HA	89.94	0.49
1:2:21:ARG:O	1:2:24:GLN:HB3	2.13	0.49
1:A:39:THR:OG1	1:A:40:ARG:N	2.50	0.49
1:5:19:ILE:HG12	1:5:79:LEU:HB2	1.95	0.49
1:L:19:ILE:HG12	1:L:79:LEU:CB	2.43	0.49
1:E:39:THR:OG1	1:E:40:ARG:N	2.45	0.49
1:E:122:ILE:HD12	1:E:122:ILE:N	2.20	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:S:154:ILE:HD13	1:T:67:LEU:HD21	1.93	0.49
1:T:64:THR:HB	1:T:76:VAL:HG11	1.94	0.49
1:C:64:THR:HB	1:C:76:VAL:HG11	1.96	0.49
1:Q:5:LYS:HE2	1:R:21:ARG:NH2	2.26	0.49
1:M:5:LYS:HE3	1:N:53:TRP:CD1	2.48	0.49
1:O:2:ASN:HD22	1:O:2:ASN:H	1.60	0.49
1:R:39:THR:OG1	1:R:40:ARG:N	2.45	0.49
1:P:19:ILE:HG12	1:P:79:LEU:HB2	1.94	0.49
1:2:54:VAL:HB	1:2:55:PRO:HD2	1.93	0.49
1:2:153:ALA:O	1:2:154:ILE:HG12	2.13	0.49
1:F:54:VAL:HB	1:F:55:PRO:HD2	1.93	0.49
1:K:153:ALA:O	1:K:154:ILE:HG12	2.13	0.49
1:9:153:ALA:O	1:9:154:ILE:HG12	2.13	0.49
1:6:56:GLY:O	1:6:59:GLU:HG2	2.13	0.49
1:J:22:PHE:HE1	1:9:131:THR:O	80.55	0.49
1:K:143:THR:HG22	1:L:55:PRO:CB	2.59	0.49
1:W:2:ASN:HD22	1:W:2:ASN:H	1.61	0.49
1:K:2:ASN:HD22	1:K:2:ASN:H	1.61	0.49
1:P:143:THR:HG22	1:Q:55:PRO:HB2	2.00	0.49
1:6:19:ILE:HG12	1:6:79:LEU:HB2	1.94	0.49
1:W:153:ALA:O	1:W:154:ILE:HG12	2.14	0.49
1:A:55:PRO:HG2	1:A:59:GLU:HG3	1.95	0.49
1:8:31:LEU:O	1:8:35:VAL:HG23	2.12	0.49
1:S:5:LYS:HE2	1:T:21:ARG:NH2	2.18	0.49
1:I:36:ASP:OD2	1:I:40:ARG:NH1	2.46	0.49
1:U:151:LEU:O	1:U:154:ILE:HG13	2.12	0.49
1:G:19:ILE:HG12	1:G:79:LEU:CB	2.43	0.49
1:R:55:PRO:HG2	1:R:59:GLU:HG3	1.99	0.49
1:D:131:THR:O	1:F:22:PHE:HE1	2.16	0.48
1:Q:23:ASN:HA	1:6:131:THR:HA	86.62	0.48
1:B:40:ARG:O	1:K:7:ASN:ND2	119.88	0.48
1:L:111:PRO:HG3	1:L:151:LEU:HD12	1.95	0.48
1:V:19:ILE:HG21	1:V:53:TRP:CZ3	2.47	0.48
1:5:67:LEU:HD21	1:9:154:ILE:HD13	1.95	0.48
1:6:2:ASN:HD21	1:7:50:THR:HG22	1.78	0.48
1:7:2:ASN:HD21	1:8:50:THR:HG22	1.78	0.48
1:U:67:LEU:HD21	1:Y:154:ILE:HD13	1.99	0.48
1:1:54:VAL:HB	1:1:55:PRO:HD2	1.95	0.48
1:N:119:THR:HG22	1:O:87:THR:HG21	2.17	0.48
1:C:141:ALA:O	1:C:144:ALA:HB3	2.24	0.48
1:E:85:GLY:HA3	2:E:157:SO4:O4	2.12	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:I:64:THR:HB	1:I:76:VAL:HG11	1.94	0.48
1:J:122:ILE:HD12	1:J:122:ILE:N	2.17	0.48
1:N:122:ILE:CD1	1:N:122:ILE:H	2.16	0.48
1:5:2:ASN:H	1:5:2:ASN:HD22	1.61	0.48
1:L:154:ILE:HD13	1:M:67:LEU:CD2	2.54	0.48
1:O:55:PRO:HG2	1:O:59:GLU:HG3	1.95	0.48
1:A:154:ILE:HD13	1:B:67:LEU:HD21	1.94	0.48
1:C:55:PRO:HG2	1:C:59:GLU:HG3	1.95	0.48
1:S:151:LEU:HD11	1:T:62:LEU:HD13	1.99	0.48
1:T:122:ILE:HD12	1:T:122:ILE:N	2.22	0.48
1:B:23:ASN:HA	1:V:131:THR:HA	18.76	0.48
1:U:19:ILE:HG12	1:U:79:LEU:CB	2.42	0.48
1:O:19:ILE:HG21	1:O:53:TRP:CZ3	2.50	0.48
1:N:5:LYS:HE3	1:O:53:TRP:CD1	2.59	0.48
1:B:41:ILE:HA	1:K:7:ASN:ND2	118.63	0.48
1:G:7:ASN:HD22	1:P:41:ILE:HD13	49.07	0.48
1:Q:19:ILE:HG12	1:Q:79:LEU:CB	2.43	0.48
1:U:41:ILE:HD13	1:V:7:ASN:HD22	48.43	0.48
1:H:153:ALA:O	1:H:154:ILE:HG12	2.12	0.48
1:X:9:ALA:HA	1:X:42:GLY:HA2	1.95	0.48
1:Y:56:GLY:O	1:Y:59:GLU:HG2	2.14	0.48
1:L:118:THR:O	1:M:87:THR:HB	2.14	0.48
1:F:119:THR:HG22	1:G:87:THR:HG21	1.95	0.48
1:5:128:ARG:NH2	2:6:157:SO4:O4	2.41	0.48
1:M:41:ILE:HA	1:7:7:ASN:ND2	96.78	0.48
1:P:56:GLY:O	1:P:59:GLU:HG2	2.13	0.48
1:Y:39:THR:OG1	1:Y:40:ARG:N	2.51	0.48
1:Q:2:ASN:HD21	1:R:50:THR:HG22	1.82	0.48
1:P:122:ILE:H	1:P:122:ILE:CD1	2.12	0.48
1:T:27:ASN:HA	1:T:30:LEU:HB2	2.03	0.48
1:B:131:THR:HA	1:V:23:ASN:HA	1.95	0.48
1:L:55:PRO:HG2	1:L:59:GLU:HG3	1.94	0.48
1:U:9:ALA:HA	1:U:42:GLY:HA2	1.94	0.48
1:R:7:ASN:ND2	1:W:41:ILE:HA	95.91	0.48
1:9:19:ILE:HG12	1:9:79:LEU:HB2	1.96	0.48
1:J:9:ALA:HA	1:J:42:GLY:HA2	2.02	0.48
1:L:1:MET:HA	1:M:46:ASP:O	2.29	0.48
1:6:31:LEU:O	1:6:32:ASP:C	2.51	0.48
1:H:64:THR:HB	1:H:76:VAL:HG11	1.96	0.48
1:Z:122:ILE:N	1:Z:122:ILE:HD12	2.14	0.48
1:5:122:ILE:HD12	1:5:122:ILE:N	2.15	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:4:64:THR:HB	1:4:76:VAL:HG11	1.96	0.48
1:B:19:ILE:HG12	1:B:79:LEU:CB	2.44	0.48
1:Y:2:ASN:H	1:Y:2:ASN:HD22	1.63	0.48
1:U:7:ASN:ND2	1:U:8:VAL:H	2.28	0.48
1:U:54:VAL:HG11	1:U:63:ALA:HB2	1.95	0.48
1:V:111:PRO:HG3	1:V:151:LEU:HD12	2.01	0.48
1:I:19:ILE:HG12	1:I:79:LEU:CB	2.44	0.48
1:9:19:ILE:HG21	1:9:53:TRP:CZ3	2.48	0.48
1:7:151:LEU:HD11	1:8:62:LEU:HD13	1.94	0.48
1:7:54:VAL:HB	1:7:55:PRO:HD2	1.94	0.48
1:W:151:LEU:HD11	1:X:62:LEU:HD13	1.96	0.48
1:N:118:THR:O	1:O:87:THR:HB	2.44	0.48
1:M:21:ARG:O	1:M:24:GLN:HB3	2.17	0.48
1:L:39:THR:OG1	1:L:40:ARG:N	2.46	0.48
1:X:54:VAL:HB	1:X:55:PRO:HD2	1.96	0.48
1:H:143:THR:HG22	1:I:55:PRO:HB2	1.94	0.48
1:3:19:ILE:HG12	1:3:79:LEU:CB	2.44	0.48
1:J:19:ILE:HG12	1:J:79:LEU:CB	2.43	0.48
1:W:31:LEU:O	1:W:35:VAL:HG23	2.14	0.48
1:D:39:THR:OG1	1:D:40:ARG:N	2.46	0.48
1:N:39:THR:OG1	1:N:40:ARG:N	2.47	0.48
1:W:5:LYS:HE3	1:X:53:TRP:CD1	2.51	0.48
1:I:21:ARG:O	1:I:24:GLN:HB3	2.15	0.48
1:E:19:ILE:HG12	1:E:79:LEU:CB	2.47	0.48
1:Z:2:ASN:HD22	1:Z:2:ASN:H	1.59	0.48
1:5:54:VAL:HB	1:5:55:PRO:HD2	1.96	0.48
1:Z:153:ALA:O	1:Z:154:ILE:HG12	2.13	0.48
1:W:21:ARG:O	1:W:24:GLN:HB3	2.20	0.48
1:L:5:LYS:HE3	1:M:53:TRP:CD1	2.62	0.48
1:A:5:LYS:HE2	1:B:21:ARG:NH2	2.24	0.48
1:Q:30:LEU:HD12	1:Q:30:LEU:HA	1.73	0.48
1:X:19:ILE:HG12	1:X:79:LEU:CB	2.45	0.48
1:K:19:ILE:HG12	1:K:79:LEU:CB	2.49	0.48
1:Y:153:ALA:O	1:Y:154:ILE:HG12	2.18	0.48
1:Z:87:THR:HG21	1:4:119:THR:HG22	1.95	0.48
1:P:39:THR:OG1	1:P:40:ARG:N	2.49	0.48
1:N:1:MET:HB2	1:O:46:ASP:O	2.19	0.48
1:G:25:PHE:CZ	1:G:126:ILE:HG23	6.11	0.48
1:W:19:ILE:HG12	1:W:79:LEU:CB	2.44	0.48
1:M:54:VAL:HG11	1:M:63:ALA:HB2	2.00	0.48
1:E:55:PRO:HG2	1:E:59:GLU:HG3	2.02	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:31:LEU:O	1:H:32:ASP:C	2.57	0.48
1:6:21:ARG:O	1:6:24:GLN:HB3	2.13	0.47
1:I:30:LEU:HA	1:I:30:LEU:HD12	1.78	0.47
1:W:54:VAL:HG11	1:W:63:ALA:HB2	1.99	0.47
1:K:54:VAL:HG11	1:K:63:ALA:HB2	1.96	0.47
1:W:67:LEU:O	1:W:70:SER:HB3	2.33	0.47
1:H:26:ILE:HD13	1:H:125:ALA:HB1	1.96	0.47
1:1:16:ALA:HB1	1:1:67:LEU:HD13	1.96	0.47
1:X:64:THR:HB	1:X:76:VAL:HG11	1.97	0.47
1:M:19:ILE:HG12	1:M:79:LEU:CB	2.44	0.47
1:R:30:LEU:O	1:R:137:GLY:HA3	2.14	0.47
1:Y:31:LEU:O	1:Y:35:VAL:HG23	2.14	0.47
1:O:19:ILE:HG12	1:O:79:LEU:HB2	1.95	0.47
1:S:19:ILE:HG12	1:S:79:LEU:CB	2.49	0.47
1:C:19:ILE:HG12	1:C:79:LEU:CB	2.44	0.47
1:B:151:LEU:HD11	1:C:62:LEU:HD13	1.96	0.47
1:3:31:LEU:O	1:3:35:VAL:HG23	2.13	0.47
1:F:87:THR:HG21	1:J:119:THR:HG22	1.95	0.47
1:U:141:ALA:O	1:U:144:ALA:HB3	2.14	0.47
1:2:42:GLY:HA3	1:2:145:LEU:HD13	1.95	0.47
1:M:131:THR:O	1:X:22:PHE:HE1	75.00	0.47
1:D:30:LEU:HD12	1:D:30:LEU:HA	1.65	0.47
1:F:31:LEU:O	1:F:35:VAL:HG23	2.14	0.47
1:C:30:LEU:HA	1:C:30:LEU:HD12	1.71	0.47
1:5:21:ARG:O	1:5:24:GLN:HB3	2.14	0.47
1:L:54:VAL:HB	1:L:55:PRO:HD2	1.96	0.47
1:C:143:THR:HG22	1:D:55:PRO:HB2	2.04	0.47
1:F:143:THR:HG22	1:G:55:PRO:HB2	1.97	0.47
1:7:19:ILE:HG12	1:7:79:LEU:HB2	1.96	0.47
1:Y:151:LEU:O	1:Y:154:ILE:HG13	2.19	0.47
1:K:9:ALA:HA	1:K:42:GLY:HA2	2.02	0.47
1:A:151:LEU:HD11	1:B:62:LEU:HD13	1.96	0.47
1:8:153:ALA:O	1:8:154:ILE:HG12	2.14	0.47
1:U:87:THR:HB	1:Y:118:THR:O	2.36	0.47
1:L:31:LEU:O	1:L:35:VAL:HG23	2.15	0.47
1:7:64:THR:HB	1:7:76:VAL:HG11	1.97	0.47
1:M:22:PHE:HD1	1:M:23:ASN:N	2.12	0.47
1:R:31:LEU:O	1:R:35:VAL:HG23	2.14	0.47
1:1:19:ILE:HG12	1:1:79:LEU:CB	2.44	0.47
1:8:30:LEU:O	1:8:137:GLY:HA3	2.14	0.47
1:5:19:ILE:HG12	1:5:79:LEU:CB	2.44	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:111:PRO:HG3	1:B:151:LEU:HD12	1.94	0.47
1:C:122:ILE:HD12	1:C:122:ILE:N	2.21	0.47
1:L:27:ASN:HA	1:L:30:LEU:HB2	1.95	0.47
1:Q:9:ALA:HA	1:Q:42:GLY:HA2	1.96	0.47
1:E:19:ILE:HG21	1:E:53:TRP:CZ3	2.49	0.47
1:V:153:ALA:O	1:V:154:ILE:HG12	2.13	0.47
1:U:41:ILE:HD13	1:V:7:ASN:ND2	48.21	0.47
1:S:55:PRO:HG2	1:S:59:GLU:HG3	1.98	0.47
1:Z:56:GLY:O	1:Z:59:GLU:HG2	2.14	0.47
1:U:39:THR:OG1	1:U:40:ARG:N	2.50	0.47
1:1:26:ILE:HD13	1:1:125:ALA:HB1	1.95	0.47
1:8:64:THR:HB	1:8:76:VAL:HG11	1.96	0.47
1:4:30:LEU:HD12	1:4:30:LEU:HA	1.78	0.47
1:T:2:ASN:H	1:T:2:ASN:HD22	1.63	0.47
1:P:46:ASP:O	1:T:1:MET:CB	2.61	0.47
1:D:143:THR:HG22	1:E:55:PRO:HB2	1.97	0.47
1:F:56:GLY:O	1:F:59:GLU:HG2	2.14	0.47
1:N:64:THR:HB	1:N:76:VAL:HG11	1.96	0.47
1:P:18:THR:HG21	1:P:64:THR:HG22	1.96	0.47
1:1:22:PHE:HD1	1:1:23:ASN:N	2.13	0.47
1:G:25:PHE:HB3	1:G:131:THR:HG23	11.98	0.47
1:A:122:ILE:HD12	1:A:122:ILE:N	2.20	0.47
1:E:30:LEU:HA	1:E:30:LEU:HD12	1.73	0.47
1:F:19:ILE:HG21	1:F:53:TRP:CZ3	2.49	0.47
1:T:30:LEU:HA	1:T:30:LEU:HD12	1.73	0.47
1:N:22:PHE:HE1	1:W:131:THR:O	71.10	0.47
1:I:2:ASN:HD21	1:J:50:THR:CG2	2.28	0.47
1:Z:19:ILE:HG12	1:Z:79:LEU:CB	2.45	0.47
1:5:19:ILE:HG21	1:5:53:TRP:CZ3	2.49	0.47
1:B:67:LEU:O	1:B:70:SER:HB3	2.23	0.47
1:S:111:PRO:HG3	1:S:151:LEU:HD12	2.05	0.47
1:M:26:ILE:HD13	1:M:125:ALA:HB1	2.06	0.47
1:G:153:ALA:O	1:G:154:ILE:HG12	2.22	0.47
1:H:39:THR:OG1	1:H:40:ARG:N	2.47	0.47
1:I:41:ILE:HD13	1:O:7:ASN:HD22	1.80	0.47
1:O:54:VAL:HG11	1:O:63:ALA:HB2	1.98	0.47
1:K:67:LEU:CD2	1:O:154:ILE:HD13	2.50	0.47
1:J:19:ILE:HG21	1:J:53:TRP:CZ3	2.57	0.47
1:M:153:ALA:O	1:M:154:ILE:HG12	2.14	0.47
1:7:30:LEU:O	1:7:137:GLY:HA3	2.14	0.47
1:9:122:ILE:N	1:9:122:ILE:HD12	2.17	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:L:5:LYS:HE2	1:M:21:ARG:NH2	2.22	0.47
1:L:5:LYS:CE	1:M:21:ARG:HH22	2.20	0.47
1:9:24:GLN:HA	1:9:27:ASN:HB2	1.97	0.47
1:M:5:LYS:CE	1:N:21:ARG:HH22	2.43	0.47
1:C:7:ASN:ND2	1:R:41:ILE:HA	2.29	0.47
1:P:87:THR:HB	1:T:119:THR:HA	1.97	0.47
1:U:55:PRO:HG2	1:U:59:GLU:HG3	1.97	0.47
1:L:19:ILE:HG21	1:L:53:TRP:CZ3	2.49	0.47
1:2:19:ILE:HG12	1:2:79:LEU:HB2	1.95	0.47
1:8:54:VAL:HB	1:8:55:PRO:HD2	1.97	0.47
1:4:34:ALA:HA	1:4:138:ALA:HA	1.97	0.47
1:A:87:THR:HG21	1:E:119:THR:HG22	1.98	0.47
1:E:30:LEU:O	1:E:137:GLY:HA3	2.17	0.47
1:6:5:LYS:HE2	1:7:21:ARG:NH2	2.26	0.47
1:Y:19:ILE:HG21	1:Y:53:TRP:CZ3	2.50	0.47
1:H:22:PHE:HD1	1:H:23:ASN:N	2.13	0.47
1:P:111:PRO:HG3	1:P:151:LEU:HD12	1.97	0.47
1:7:55:PRO:HG2	1:7:59:GLU:HG3	1.96	0.47
1:W:26:ILE:HD13	1:W:125:ALA:HB1	2.02	0.47
1:F:18:THR:HG21	1:F:64:THR:HG22	2.04	0.46
1:Q:131:THR:HA	1:1:23:ASN:OD1	86.18	0.46
1:1:122:ILE:HD12	1:1:122:ILE:N	2.17	0.46
1:4:122:ILE:HD12	1:4:122:ILE:N	2.20	0.46
1:G:5:LYS:CE	1:H:21:ARG:HH22	2.23	0.46
1:Z:5:LYS:CE	1:1:21:ARG:HH22	2.22	0.46
1:M:27:ASN:HA	1:M:30:LEU:HB2	1.97	0.46
1:9:30:LEU:O	1:9:137:GLY:HA3	2.15	0.46
1:B:21:ARG:O	1:B:24:GLN:HB3	2.14	0.46
1:N:27:ASN:HA	1:N:30:LEU:HB2	1.97	0.46
1:K:27:ASN:HA	1:K:30:LEU:HB2	1.98	0.46
1:N:22:PHE:HD1	1:N:23:ASN:N	2.15	0.46
1:P:23:ASN:HA	1:3:130:GLY:O	2.14	0.46
1:N:54:VAL:HG11	1:N:63:ALA:HB2	1.96	0.46
1:E:54:VAL:HB	1:E:55:PRO:HD2	2.00	0.46
1:8:27:ASN:HA	1:8:30:LEU:HB2	1.97	0.46
1:6:70:SER:OG	1:6:72:LYS:HB2	2.14	0.46
1:U:46:ASP:O	1:Y:1:MET:HA	2.15	0.46
1:Z:18:THR:HG21	1:Z:64:THR:HG22	1.97	0.46
1:M:131:THR:O	1:8:22:PHE:HE1	71.58	0.46
1:F:27:ASN:HA	1:F:30:LEU:HB2	1.97	0.46
1:T:21:ARG:O	1:T:24:GLN:HB3	2.15	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:W:1:MET:CB	1:X:46:ASP:O	2.98	0.46
1:A:50:THR:CG2	1:E:2:ASN:HD21	2.25	0.46
1:I:31:LEU:O	1:I:35:VAL:HG23	2.19	0.46
1:8:19:ILE:HD12	1:8:31:LEU:HD22	1.95	0.46
1:9:31:LEU:O	1:9:35:VAL:HG23	2.15	0.46
1:3:143:THR:HG22	1:4:55:PRO:HB2	1.97	0.46
1:4:55:PRO:HG2	1:4:59:GLU:HG3	1.97	0.46
1:T:30:LEU:O	1:T:137:GLY:HA3	2.20	0.46
1:P:31:LEU:O	1:P:35:VAL:HG23	2.17	0.46
1:J:23:ASN:OD1	1:9:131:THR:HA	79.65	0.46
1:B:9:ALA:HA	1:B:42:GLY:HA2	1.98	0.46
1:T:39:THR:OG1	1:T:40:ARG:N	2.49	0.46
1:G:7:ASN:ND2	1:G:8:VAL:H	2.13	0.46
1:R:54:VAL:HB	1:R:55:PRO:HD2	2.02	0.46
1:S:118:THR:O	1:T:87:THR:HB	2.34	0.46
1:4:19:ILE:HG12	1:4:79:LEU:CB	2.46	0.46
1:1:55:PRO:HG2	1:1:59:GLU:HG3	1.96	0.46
1:G:31:LEU:O	1:G:35:VAL:HG23	2.16	0.46
1:W:22:PHE:HD1	1:W:23:ASN:N	2.13	0.46
1:Y:22:PHE:HD1	1:Y:23:ASN:N	2.17	0.46
1:I:111:PRO:HG3	1:I:151:LEU:HD12	1.99	0.46
1:Y:54:VAL:HB	1:Y:55:PRO:HD2	1.98	0.46
1:X:62:LEU:HD23	1:X:62:LEU:HA	1.80	0.46
1:P:64:THR:HB	1:P:76:VAL:HG11	2.00	0.46
1:F:151:LEU:O	1:F:154:ILE:HG13	2.21	0.46
1:G:9:ALA:HA	1:G:42:GLY:HA2	1.97	0.46
1:V:9:ALA:HA	1:V:42:GLY:HA2	1.99	0.46
1:2:111:PRO:HG3	1:2:151:LEU:HD12	1.96	0.46
1:D:70:SER:OG	1:D:72:LYS:HB2	2.15	0.46
1:B:1:MET:HB2	1:C:46:ASP:O	2.15	0.46
1:S:30:LEU:HD12	1:S:30:LEU:HA	1.75	0.46
1:T:54:VAL:HG11	1:T:63:ALA:HB2	1.98	0.46
1:V:54:VAL:HG11	1:V:63:ALA:HB2	2.01	0.46
1:C:41:ILE:HA	1:R:7:ASN:ND2	2.31	0.46
1:3:9:ALA:HA	1:3:42:GLY:HA2	1.98	0.46
1:Z:87:THR:HB	1:4:118:THR:O	2.16	0.46
2:B:159:SO4:O2	1:C:85:GLY:HA3	2.16	0.46
1:G:18:THR:HG21	1:G:64:THR:HG22	1.97	0.46
1:Y:122:ILE:HD12	1:Y:122:ILE:N	2.18	0.46
1:H:127:GLU:OE1	2:H:158:SO4:O2	2.66	0.46
1:O:27:ASN:HA	1:O:30:LEU:HB2	1.99	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:Q:22:PHE:HD1	1:Q:23:ASN:N	2.14	0.46
1:V:7:ASN:OD1	1:V:9:ALA:HB3	2.15	0.46
1:2:151:LEU:HD11	1:3:62:LEU:HD13	1.98	0.46
1:8:19:ILE:HG12	1:8:79:LEU:CB	2.46	0.46
1:G:122:ILE:HD13	1:G:123:GLU:OE2	9.58	0.46
1:C:27:ASN:HA	1:C:30:LEU:HB2	1.96	0.46
1:X:30:LEU:O	1:X:137:GLY:HA3	2.16	0.46
1:V:22:PHE:HD1	1:V:23:ASN:N	2.13	0.46
1:C:22:PHE:HE1	1:K:131:THR:O	87.90	0.46
1:A:22:PHE:HD1	1:A:23:ASN:N	2.14	0.46
1:K:54:VAL:HB	1:K:55:PRO:HD2	1.98	0.46
1:P:154:ILE:HD13	1:Q:67:LEU:HD21	2.01	0.46
1:M:111:PRO:HG3	1:M:151:LEU:HD12	2.02	0.46
1:C:151:LEU:O	1:C:154:ILE:HG13	2.15	0.46
1:M:1:MET:HB2	1:N:46:ASP:O	2.16	0.46
1:K:111:PRO:HG3	1:K:151:LEU:HD12	1.99	0.46
1:S:154:ILE:HD13	1:T:67:LEU:CD2	2.48	0.46
1:5:27:ASN:HA	1:5:30:LEU:HB2	1.98	0.46
1:U:30:LEU:HA	1:U:30:LEU:HD12	1.71	0.46
1:U:30:LEU:O	1:U:137:GLY:HA3	2.15	0.46
1:Q:151:LEU:O	1:Q:154:ILE:HG13	2.25	0.46
1:B:151:LEU:O	1:B:154:ILE:HG13	2.15	0.46
1:5:153:ALA:O	1:5:154:ILE:HG12	2.16	0.46
1:3:54:VAL:HG11	1:3:63:ALA:HB2	1.98	0.46
1:F:53:TRP:CD1	1:J:5:LYS:HE3	2.61	0.46
1:W:9:ALA:HA	1:W:42:GLY:HA2	2.03	0.46
1:Y:27:ASN:HA	1:Y:30:LEU:HB2	1.98	0.46
1:X:54:VAL:HG11	1:X:63:ALA:HB2	2.00	0.46
1:N:9:ALA:HA	1:N:42:GLY:HA2	2.00	0.46
1:O:9:ALA:HA	1:O:42:GLY:HA2	1.98	0.46
1:2:154:ILE:HD13	1:3:67:LEU:HD21	1.98	0.46
1:F:7:ASN:ND2	1:G:41:ILE:HA	35.62	0.46
1:B:54:VAL:HG11	1:B:63:ALA:HB2	1.99	0.46
1:3:153:ALA:O	1:3:154:ILE:HG12	2.16	0.46
1:A:153:ALA:O	1:A:154:ILE:HG12	2.16	0.46
1:X:111:PRO:HG3	1:X:151:LEU:HD12	1.98	0.46
1:S:41:ILE:HA	1:Y:7:ASN:ND2	2.31	0.46
1:7:31:LEU:O	1:7:35:VAL:HG23	2.16	0.46
1:P:26:ILE:HD13	1:P:125:ALA:HB1	1.98	0.46
1:J:21:ARG:O	1:J:24:GLN:HB3	2.17	0.45
1:D:154:ILE:HD13	1:E:67:LEU:HD21	1.98	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:18:THR:HG21	1:H:64:THR:HG22	2.02	0.45
1:C:7:ASN:OD1	1:C:9:ALA:HB3	2.18	0.45
1:L:22:PHE:HD1	1:L:23:ASN:N	2.16	0.45
1:G:56:GLY:O	1:G:59:GLU:HG2	2.16	0.45
1:P:62:LEU:HA	1:P:62:LEU:HD23	1.75	0.45
1:1:56:GLY:O	1:1:59:GLU:HG2	2.15	0.45
1:A:118:THR:O	1:B:87:THR:HB	2.17	0.45
1:F:141:ALA:O	1:F:144:ALA:HB3	2.16	0.45
1:F:30:LEU:O	1:F:137:GLY:HA3	2.16	0.45
1:X:27:ASN:HA	1:X:30:LEU:HB2	1.98	0.45
1:K:119:THR:HG22	1:L:87:THR:HG21	2.05	0.45
1:R:19:ILE:HG12	1:R:79:LEU:CB	2.49	0.45
1:2:56:GLY:O	1:2:59:GLU:HG2	2.17	0.45
1:E:41:ILE:HA	1:8:7:ASN:ND2	120.38	0.45
1:S:31:LEU:O	1:S:34:ALA:N	2.57	0.45
1:C:56:GLY:O	1:C:59:GLU:HG2	2.17	0.45
1:1:1:MET:HB2	1:2:46:ASP:O	2.15	0.45
1:F:26:ILE:HD13	1:F:125:ALA:HB1	2.02	0.45
1:8:119:THR:HG22	1:9:87:THR:HG21	1.97	0.45
1:7:122:ILE:HD12	1:7:122:ILE:N	2.17	0.45
1:U:122:ILE:HD12	1:U:122:ILE:N	2.19	0.45
1:U:126:ILE:H	1:U:126:ILE:HG12	1.63	0.45
1:R:37:ALA:O	1:R:41:ILE:HB	2.17	0.45
1:F:111:PRO:HG3	1:F:151:LEU:HD12	1.99	0.45
1:G:67:LEU:O	1:G:70:SER:HB3	2.30	0.45
1:N:54:VAL:HB	1:N:55:PRO:HD2	1.99	0.45
1:R:143:THR:HG22	1:S:55:PRO:HB2	2.02	0.45
1:P:67:LEU:O	1:P:70:SER:HB3	2.23	0.45
1:F:7:ASN:OD1	1:F:9:ALA:HB3	2.23	0.45
1:7:154:ILE:HD13	1:8:67:LEU:HD21	1.97	0.45
1:Y:50:THR:CG2	1:Y:73:TYR:HE1	2.34	0.45
1:U:1:MET:HB2	1:V:46:ASP:O	2.16	0.45
1:7:143:THR:HG22	1:8:55:PRO:HB2	1.99	0.45
1:8:56:GLY:O	1:8:59:GLU:HG2	2.16	0.45
1:G:22:PHE:HD1	1:G:23:ASN:N	2.13	0.45
1:X:19:ILE:HD12	1:X:31:LEU:HD22	1.98	0.45
1:P:19:ILE:HG12	1:P:79:LEU:CB	2.46	0.45
1:S:22:PHE:HD1	1:S:23:ASN:N	2.15	0.45
1:J:31:LEU:O	1:J:35:VAL:HG23	2.17	0.45
1:9:54:VAL:HG11	1:9:63:ALA:HB2	1.97	0.45
1:U:62:LEU:HA	1:U:62:LEU:HD23	1.87	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:1:39:THR:OG1	1:1:40:ARG:N	2.50	0.45
1:Q:130:GLY:O	1:1:23:ASN:HA	89.02	0.45
1:V:30:LEU:O	1:V:137:GLY:HA3	2.22	0.45
1:N:19:ILE:HD12	1:N:31:LEU:HD22	2.09	0.45
1:O:30:LEU:O	1:O:137:GLY:HA3	2.24	0.45
1:P:54:VAL:HB	1:P:55:PRO:HD2	1.99	0.45
1:O:7:ASN:ND2	1:O:8:VAL:H	2.25	0.45
1:H:111:PRO:HG3	1:H:151:LEU:HD12	1.99	0.45
1:S:9:ALA:HA	1:S:42:GLY:HA2	1.97	0.45
1:J:54:VAL:HG11	1:J:63:ALA:HB2	2.00	0.45
1:5:50:THR:HG22	1:9:2:ASN:HD21	1.81	0.45
1:H:7:ASN:ND2	1:2:40:ARG:O	100.42	0.45
1:H:31:LEU:O	1:H:35:VAL:HG23	2.16	0.45
1:F:36:ASP:OD2	1:F:40:ARG:NH1	2.60	0.45
1:S:39:THR:OG1	1:S:40:ARG:N	2.51	0.45
1:F:1:MET:HB2	1:G:46:ASP:O	2.16	0.45
1:M:18:THR:HG21	1:M:64:THR:HG22	1.98	0.45
1:D:27:ASN:HA	1:D:30:LEU:HB2	1.97	0.45
1:N:126:ILE:HG12	1:N:126:ILE:H	1.68	0.45
1:B:124:GLN:NE2	2:B:158:SO4:O2	2.65	0.45
1:C:30:LEU:O	1:C:137:GLY:HA3	2.16	0.45
1:M:41:ILE:HA	1:W:7:ASN:ND2	99.94	0.45
1:G:30:LEU:HA	1:G:30:LEU:HD12	1.73	0.45
1:C:22:PHE:HD1	1:C:23:ASN:N	2.18	0.45
1:P:23:ASN:HA	1:3:131:THR:HA	1.98	0.45
1:L:7:ASN:OD1	1:L:9:ALA:HB3	2.17	0.45
1:F:54:VAL:HG11	1:F:63:ALA:HB2	1.98	0.45
1:L:67:LEU:O	1:L:70:SER:HB3	2.16	0.45
1:A:19:ILE:HG12	1:A:79:LEU:CB	2.49	0.45
1:1:111:PRO:HG3	1:1:151:LEU:HD12	1.99	0.45
1:K:1:MET:HA	1:L:46:ASP:O	2.28	0.45
1:F:62:LEU:HD13	1:J:151:LEU:HD11	1.98	0.45
1:9:11:PRO:HB3	1:9:43:GLN:HB3	1.97	0.45
1:2:122:ILE:HD12	1:2:122:ILE:N	2.19	0.45
1:2:5:LYS:CE	1:3:21:ARG:HH22	2.21	0.45
1:G:5:LYS:HE3	1:H:53:TRP:CD1	2.59	0.45
1:M:31:LEU:O	1:M:32:ASP:C	2.66	0.45
1:A:5:LYS:HE3	1:B:53:TRP:CD1	2.67	0.45
1:E:131:THR:O	1:O:22:PHE:HE1	1.99	0.45
1:Y:31:LEU:O	1:Y:34:ALA:N	2.63	0.45
1:3:22:PHE:HD1	1:3:23:ASN:N	2.15	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:Y:9:ALA:HA	1:Y:42:GLY:HA2	1.97	0.45
1:P:20:ALA:HA	1:P:54:VAL:O	2.16	0.45
1:X:131:THR:HA	1:Z:23:ASN:HA	1.99	0.45
1:E:7:ASN:OD1	1:E:9:ALA:HB3	2.24	0.45
1:A:40:ARG:O	1:L:7:ASN:ND2	73.59	0.45
1:H:54:VAL:HG11	1:H:63:ALA:HB2	1.99	0.45
1:I:7:ASN:ND2	1:9:41:ILE:HA	110.53	0.45
1:2:19:ILE:HG12	1:2:79:LEU:CB	2.47	0.45
1:D:30:LEU:O	1:D:137:GLY:HA3	2.21	0.45
1:S:27:ASN:HA	1:S:30:LEU:HB2	1.98	0.45
1:N:30:LEU:O	1:N:137:GLY:HA3	2.20	0.45
1:G:27:ASN:HA	1:G:30:LEU:HB2	2.01	0.45
1:Z:27:ASN:HA	1:Z:30:LEU:HB2	1.99	0.45
1:O:19:ILE:HD11	1:O:31:LEU:HB2	1.98	0.45
1:G:151:LEU:O	1:G:154:ILE:HG13	2.19	0.45
1:S:23:ASN:HA	1:W:130:GLY:O	70.62	0.45
1:D:9:ALA:HA	1:D:42:GLY:HA2	2.03	0.45
1:U:87:THR:HG21	1:Y:119:THR:HG22	2.09	0.45
1:B:128:ARG:NH2	2:B:159:SO4:O1	4.63	0.45
1:K:151:LEU:HD11	1:L:62:LEU:HD13	1.99	0.45
1:Z:151:LEU:HD11	1:1:62:LEU:HD13	1.99	0.45
1:W:39:THR:OG1	1:W:40:ARG:N	2.50	0.45
1:D:119:THR:HG22	1:E:87:THR:HG21	1.98	0.45
1:N:154:ILE:HD13	1:O:67:LEU:HD21	1.99	0.45
1:B:30:LEU:HD12	1:B:30:LEU:HA	1.73	0.45
1:C:7:ASN:ND2	1:R:40:ARG:O	2.50	0.45
1:L:30:LEU:O	1:L:137:GLY:HA3	2.17	0.45
1:P:22:PHE:HD1	1:P:23:ASN:N	2.15	0.45
1:E:54:VAL:HG11	1:E:63:ALA:HB2	2.00	0.45
1:P:7:ASN:OD1	1:P:9:ALA:HB3	2.17	0.45
1:5:151:LEU:HD11	1:6:62:LEU:HD13	1.98	0.45
1:M:39:THR:OG1	1:M:40:ARG:N	2.50	0.45
1:7:124:GLN:HG2	2:7:158:SO4:O2	2.17	0.45
1:I:5:LYS:NZ	1:J:21:ARG:NH2	2.89	0.45
1:D:151:LEU:HD11	1:E:62:LEU:HD13	1.99	0.45
1:B:22:PHE:HD1	1:B:23:ASN:N	2.15	0.45
1:F:153:ALA:C	1:F:154:ILE:HG12	2.37	0.45
1:Z:30:LEU:HA	1:Z:30:LEU:HD12	1.74	0.45
1:Y:22:PHE:HE1	1:4:131:THR:O	2.00	0.45
1:G:54:VAL:HG11	1:G:63:ALA:HB2	1.98	0.45
1:Z:19:ILE:HD12	1:Z:31:LEU:HD22	1.99	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:143:THR:HG22	1:B:55:PRO:HB2	1.99	0.45
1:5:56:GLY:O	1:5:59:GLU:HG2	2.16	0.45
1:Z:143:THR:HG22	1:1:55:PRO:HB2	1.99	0.45
1:D:118:THR:O	1:E:87:THR:HB	2.17	0.45
1:K:31:LEU:O	1:K:35:VAL:HG23	2.17	0.45
1:H:141:ALA:O	1:H:144:ALA:HB3	2.17	0.45
1:O:141:ALA:O	1:O:144:ALA:HB3	2.27	0.45
1:S:21:ARG:O	1:S:24:GLN:HB3	2.16	0.44
1:S:5:LYS:CE	1:T:21:ARG:HH22	2.20	0.44
1:N:31:LEU:O	1:N:35:VAL:HG23	2.17	0.44
1:H:40:ARG:O	1:2:7:ASN:ND2	96.34	0.44
1:K:55:PRO:HG2	1:K:59:GLU:HG3	1.99	0.44
1:H:151:LEU:HD11	1:I:62:LEU:HD13	2.08	0.44
1:H:154:ILE:HD13	1:I:67:LEU:HD21	2.10	0.44
1:2:151:LEU:O	1:2:154:ILE:HG13	2.16	0.44
1:J:55:PRO:HG2	1:J:59:GLU:HG3	2.00	0.44
1:1:9:ALA:HA	1:1:42:GLY:HA2	1.99	0.44
1:F:39:THR:OG1	1:F:40:ARG:N	2.53	0.44
1:K:26:ILE:HD13	1:K:125:ALA:HB1	2.02	0.44
1:G:34:ALA:HA	1:G:138:ALA:HA	2.02	0.44
1:Z:39:THR:OG1	1:Z:40:ARG:N	2.50	0.44
1:D:31:LEU:O	1:D:35:VAL:HG23	2.22	0.44
1:W:27:ASN:HA	1:W:30:LEU:HB2	1.99	0.44
1:3:21:ARG:O	1:3:24:GLN:HB3	2.17	0.44
1:5:30:LEU:HD12	1:5:30:LEU:HA	1.74	0.44
1:I:54:VAL:HG11	1:I:63:ALA:HB2	1.98	0.44
1:3:62:LEU:HA	1:3:62:LEU:HD23	1.78	0.44
1:6:54:VAL:HB	1:6:55:PRO:HD2	1.98	0.44
1:P:1:MET:HA	1:Q:46:ASP:O	2.17	0.44
1:J:64:THR:HB	1:J:76:VAL:HG11	1.98	0.44
1:L:122:ILE:HD13	1:L:123:GLU:OE2	9.53	0.44
1:Y:18:THR:OG1	1:Y:64:THR:HG22	2.25	0.44
1:H:30:LEU:HA	1:H:30:LEU:HD12	1.78	0.44
1:H:27:ASN:HA	1:H:30:LEU:HB2	1.98	0.44
1:P:30:LEU:O	1:P:137:GLY:HA3	2.18	0.44
1:N:7:ASN:OD1	1:N:9:ALA:HB3	2.20	0.44
1:F:9:ALA:HA	1:F:42:GLY:HA2	2.00	0.44
1:E:151:LEU:O	1:E:154:ILE:HG13	2.17	0.44
1:5:54:VAL:HG11	1:5:63:ALA:HB2	1.98	0.44
1:F:34:ALA:HA	1:F:138:ALA:HA	2.00	0.44
1:B:30:LEU:O	1:B:137:GLY:HA3	2.17	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:K:30:LEU:O	1:K:137:GLY:HA3	2.17	0.44
1:F:151:LEU:O	1:F:154:ILE:N	2.67	0.44
1:O:19:ILE:HG12	1:O:79:LEU:CB	2.48	0.44
1:7:22:PHE:HD1	1:7:23:ASN:N	2.15	0.44
1:2:54:VAL:HG11	1:2:63:ALA:HB2	2.00	0.44
1:D:7:ASN:OD1	1:D:9:ALA:HB3	2.17	0.44
1:K:1:MET:HB2	1:L:46:ASP:O	2.22	0.44
1:F:122:ILE:N	1:F:122:ILE:HD12	2.21	0.44
1:P:30:LEU:HD12	1:P:30:LEU:HA	1.81	0.44
1:P:46:ASP:O	1:T:1:MET:HA	2.23	0.44
1:M:20:ALA:HA	1:M:54:VAL:O	2.18	0.44
1:C:39:THR:OG1	1:C:40:ARG:N	2.51	0.44
1:S:7:ASN:ND2	1:S:8:VAL:H	2.20	0.44
1:T:62:LEU:HD23	1:T:62:LEU:HA	1.80	0.44
1:Z:111:PRO:HG3	1:Z:151:LEU:HD12	1.99	0.44
1:K:31:LEU:O	1:K:34:ALA:N	2.55	0.44
1:S:26:ILE:HD13	1:S:125:ALA:HB1	2.01	0.44
1:Q:1:MET:HA	1:R:46:ASP:O	2.36	0.44
1:V:141:ALA:O	1:V:144:ALA:HB3	2.17	0.44
1:F:21:ARG:NH2	1:J:5:LYS:HE2	2.22	0.44
1:P:5:LYS:CE	1:Q:21:ARG:HH22	2.22	0.44
1:T:153:ALA:C	1:T:154:ILE:HG12	2.45	0.44
1:A:37:ALA:O	1:A:41:ILE:HB	2.22	0.44
1:A:36:ASP:OD2	1:A:40:ARG:NH1	2.57	0.44
1:R:54:VAL:HG11	1:R:63:ALA:HB2	1.98	0.44
1:Y:111:PRO:HG3	1:Y:151:LEU:HD12	2.01	0.44
1:4:153:ALA:O	1:4:154:ILE:HG12	2.17	0.44
1:A:87:THR:HB	1:E:118:THR:O	2.17	0.44
1:1:64:THR:HB	1:1:76:VAL:HG11	1.99	0.44
1:R:18:THR:HG21	1:R:64:THR:HG22	2.00	0.44
1:L:18:THR:HG21	1:L:64:THR:HG22	1.99	0.44
1:H:5:LYS:HE2	1:I:21:ARG:NH2	2.40	0.44
1:5:85:GLY:HA3	2:5:157:SO4:S	2.57	0.44
1:F:22:PHE:HD1	1:F:23:ASN:N	2.17	0.44
1:R:85:GLY:HA3	2:R:157:SO4:O4	2.17	0.44
1:B:40:ARG:O	1:U:7:ASN:ND2	2.50	0.44
1:L:7:ASN:ND2	1:L:8:VAL:H	2.21	0.44
1:P:41:ILE:HA	1:6:7:ASN:ND2	110.38	0.44
1:R:151:LEU:O	1:R:154:ILE:HG13	2.18	0.44
1:R:50:THR:CG2	1:R:73:TYR:HE1	2.38	0.44
1:R:119:THR:HG22	1:S:87:THR:HG21	1.99	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:4:31:LEU:O	1:4:35:VAL:HG23	2.18	0.44
1:6:34:ALA:HA	1:6:138:ALA:HA	2.00	0.44
1:G:126:ILE:H	1:G:126:ILE:HG12	1.69	0.44
1:4:18:THR:HG21	1:4:64:THR:HG22	1.99	0.44
1:R:27:ASN:HA	1:R:30:LEU:HB2	2.03	0.44
1:O:30:LEU:HD12	1:O:30:LEU:HA	1.67	0.44
1:4:22:PHE:HD1	1:4:23:ASN:N	2.16	0.44
1:U:131:THR:O	1:W:22:PHE:HE1	40.49	0.44
1:U:142:LEU:HD12	1:U:142:LEU:HA	1.87	0.44
1:Q:54:VAL:HG11	1:Q:63:ALA:HB2	1.99	0.44
1:S:54:VAL:HG11	1:S:63:ALA:HB2	2.01	0.44
1:I:9:ALA:HA	1:I:42:GLY:HA2	2.00	0.44
1:3:154:ILE:HD13	1:4:67:LEU:HD21	2.00	0.44
1:G:36:ASP:OD2	1:G:40:ARG:NH1	2.53	0.44
2:7:158:SO4:O3	1:8:86:GLY:HA3	2.18	0.44
1:6:119:THR:HG22	1:7:87:THR:HG21	2.00	0.44
1:1:31:LEU:O	1:1:34:ALA:N	2.51	0.44
1:S:105:ALA:HB2	1:S:112:VAL:HG23	2.05	0.44
1:D:25:PHE:CD1	1:D:25:PHE:C	2.91	0.44
1:O:25:PHE:CZ	1:T:126:ILE:HG23	81.52	0.44
1:T:126:ILE:H	1:T:126:ILE:HG12	1.68	0.44
1:X:30:LEU:HD12	1:X:30:LEU:HA	1.69	0.44
1:M:50:THR:CG2	1:M:73:TYR:HE1	2.34	0.44
1:6:22:PHE:HD1	1:6:23:ASN:N	2.16	0.44
1:M:62:LEU:HD23	1:M:62:LEU:HA	1.83	0.44
1:C:40:ARG:O	1:R:7:ASN:ND2	2.50	0.44
1:8:30:LEU:HA	1:8:30:LEU:HD12	1.71	0.44
1:I:5:LYS:CE	1:J:21:ARG:NH2	2.96	0.43
1:Y:83:ILE:HD13	1:Y:122:ILE:HA	2.01	0.43
1:L:30:LEU:HA	1:L:30:LEU:HD12	1.72	0.43
1:Q:7:ASN:ND2	1:Q:8:VAL:H	2.27	0.43
1:X:50:THR:CG2	1:X:73:TYR:HE1	2.31	0.43
1:O:20:ALA:HA	1:O:54:VAL:O	2.24	0.43
1:A:30:LEU:O	1:A:137:GLY:HA3	2.21	0.43
1:A:27:ASN:HA	1:A:30:LEU:HB2	2.02	0.43
1:3:18:THR:HG21	1:3:64:THR:HG22	2.00	0.43
1:F:31:LEU:O	1:F:32:ASP:C	2.56	0.43
1:H:19:ILE:HG12	1:H:79:LEU:CB	2.49	0.43
1:5:22:PHE:HD1	1:5:23:ASN:N	2.16	0.43
1:Q:7:ASN:OD1	1:Q:9:ALA:HB3	2.22	0.43
1:D:22:PHE:HD1	1:D:23:ASN:N	2.15	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:6:19:ILE:HG12	1:6:79:LEU:CB	2.48	0.43
1:D:26:ILE:HD13	1:D:125:ALA:HB1	2.06	0.43
1:M:126:ILE:H	1:M:126:ILE:HG12	1.71	0.43
1:M:60:LEU:HA	1:M:60:LEU:HD23	1.81	0.43
1:M:5:LYS:HE2	1:N:21:ARG:NH2	2.50	0.43
1:O:31:LEU:O	1:O:32:ASP:C	2.56	0.43
1:A:7:ASN:OD1	1:A:9:ALA:HB3	2.20	0.43
1:R:154:ILE:HD13	1:S:67:LEU:HD21	2.12	0.43
1:A:111:PRO:HG3	1:A:151:LEU:HD12	2.00	0.43
1:1:7:ASN:OD1	1:1:9:ALA:HB3	2.18	0.43
1:6:54:VAL:HG11	1:6:63:ALA:HB2	1.99	0.43
1:J:111:PRO:HG3	1:J:151:LEU:HD12	2.00	0.43
1:E:26:ILE:HD13	1:E:125:ALA:HB1	2.00	0.43
1:S:123:GLU:OE2	1:U:122:ILE:HD13	2.33	0.43
1:C:23:ASN:OD1	1:K:131:THR:HA	86.22	0.43
1:J:23:ASN:HA	1:O:130:GLY:O	2.19	0.43
1:N:130:GLY:O	1:P:23:ASN:HA	92.28	0.43
1:E:22:PHE:HD1	1:E:23:ASN:N	2.15	0.43
1:P:37:ALA:O	1:P:41:ILE:HB	2.18	0.43
1:M:118:THR:HG21	1:N:92:TYR:CE2	2.76	0.43
1:Y:54:VAL:HG11	1:Y:63:ALA:HB2	2.00	0.43
1:K:154:ILE:HD13	1:L:67:LEU:CD2	2.56	0.43
1:5:55:PRO:HB2	1:9:143:THR:HG22	2.00	0.43
1:P:34:ALA:HA	1:P:138:ALA:HA	2.00	0.43
1:D:141:ALA:O	1:D:144:ALA:HB3	2.29	0.43
1:X:22:PHE:HD1	1:X:23:ASN:N	2.18	0.43
1:L:126:ILE:HG12	1:L:126:ILE:H	1.74	0.43
1:S:30:LEU:O	1:S:137:GLY:HA3	2.23	0.43
1:L:2:ASN:HD21	1:M:50:THR:CG2	2.35	0.43
1:Z:19:ILE:HG21	1:Z:53:TRP:CZ3	2.53	0.43
1:7:54:VAL:HG11	1:7:63:ALA:HB2	2.00	0.43
1:I:153:ALA:O	1:I:154:ILE:HG12	2.20	0.43
1:G:26:ILE:HD13	1:G:125:ALA:HB1	2.02	0.43
1:Y:26:ILE:HD13	1:Y:125:ALA:HB1	2.05	0.43
1:7:62:LEU:HD23	1:7:62:LEU:HA	1.75	0.43
1:2:62:LEU:HD23	1:2:62:LEU:HA	1.79	0.43
1:N:18:THR:HG21	1:N:64:THR:HG22	2.00	0.43
1:E:18:THR:HG21	1:E:64:THR:HG22	2.03	0.43
1:6:27:ASN:HA	1:6:30:LEU:HB2	2.01	0.43
1:X:31:LEU:O	1:X:35:VAL:HG23	2.19	0.43
1:K:30:LEU:HD12	1:K:30:LEU:HA	1.79	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:30:LEU:O	1:G:137:GLY:HA3	2.19	0.43
1:Y:31:LEU:O	1:Y:32:ASP:C	2.55	0.43
1:I:22:PHE:HD1	1:I:23:ASN:N	2.17	0.43
1:I:40:ARG:O	1:9:7:ASN:ND2	112.86	0.43
1:M:7:ASN:ND2	1:W:41:ILE:HA	96.50	0.43
1:8:36:ASP:OD2	1:8:40:ARG:NH1	2.52	0.43
1:J:36:ASP:OD2	1:J:40:ARG:NH1	2.51	0.43
1:G:85:GLY:HA3	2:G:157:SO4:O2	2.18	0.43
1:S:18:THR:OG1	1:S:64:THR:HG22	2.22	0.43
1:A:126:ILE:HG12	1:A:126:ILE:H	1.68	0.43
1:F:126:ILE:H	1:F:126:ILE:HG12	1.62	0.43
1:8:18:THR:HG21	1:8:64:THR:HG22	1.99	0.43
1:T:20:ALA:HA	1:T:54:VAL:O	2.21	0.43
1:Y:19:ILE:HG12	1:Y:79:LEU:CB	2.48	0.43
1:B:130:GLY:O	1:L:23:ASN:HA	88.65	0.43
1:J:22:PHE:HD1	1:J:23:ASN:N	2.16	0.43
1:J:142:LEU:HA	1:J:142:LEU:HD12	1.87	0.43
1:1:143:THR:HG22	1:2:55:PRO:CB	2.46	0.43
1:X:154:ILE:HD13	1:Y:67:LEU:HD21	2.02	0.43
1:P:151:LEU:HD11	1:Q:62:LEU:HD13	2.01	0.43
1:Y:20:ALA:HA	1:Y:54:VAL:O	2.19	0.43
1:4:7:ASN:OD1	1:4:9:ALA:HB3	2.17	0.43
1:Q:26:ILE:HD13	1:Q:125:ALA:HB1	2.00	0.43
1:J:30:LEU:O	1:J:137:GLY:HA3	2.19	0.43
1:O:83:ILE:HD13	1:O:122:ILE:HA	2.07	0.43
1:Q:18:THR:HG21	1:Q:64:THR:HG22	2.00	0.43
1:9:27:ASN:HA	1:9:30:LEU:HB2	2.01	0.43
1:R:30:LEU:HD12	1:R:30:LEU:HA	1.80	0.43
1:5:30:LEU:O	1:5:137:GLY:HA3	2.19	0.43
1:U:31:LEU:O	1:U:35:VAL:HG23	2.19	0.43
1:P:23:ASN:OD1	1:3:131:THR:HA	2.19	0.43
1:I:143:THR:HG22	1:J:55:PRO:HB2	2.00	0.43
1:S:34:ALA:HA	1:S:138:ALA:HA	2.05	0.43
1:B:31:LEU:O	1:B:34:ALA:N	2.52	0.43
1:L:62:LEU:HD23	1:L:62:LEU:HA	1.83	0.43
1:2:31:LEU:O	1:2:32:ASP:C	2.57	0.43
1:5:39:THR:OG1	1:5:40:ARG:N	2.51	0.43
1:E:31:LEU:O	1:E:35:VAL:HG23	2.23	0.43
1:8:85:GLY:HA3	2:8:157:SO4:O4	2.18	0.43
1:F:83:ILE:HD13	1:F:122:ILE:HA	2.04	0.43
1:T:16:ALA:O	1:T:76:VAL:HA	2.19	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:M:123:GLU:OE2	1:X:122:ILE:HD13	65.20	0.43
1:W:60:LEU:HD23	1:W:60:LEU:HA	1.86	0.43
1:M:30:LEU:O	1:M:137:GLY:HA3	2.21	0.43
1:K:22:PHE:HD1	1:K:23:ASN:N	2.18	0.43
1:G:111:PRO:HG3	1:G:151:LEU:HD12	2.00	0.43
1:S:131:THR:HA	1:U:23:ASN:HA	2.11	0.43
1:Z:54:VAL:HG11	1:Z:63:ALA:HB2	2.00	0.43
1:C:151:LEU:HD11	1:D:62:LEU:HD13	2.04	0.43
1:N:36:ASP:OD2	1:N:40:ARG:NH1	2.64	0.43
1:O:105:ALA:HB2	1:O:112:VAL:HG23	2.11	0.43
1:3:151:LEU:HD11	1:4:62:LEU:HD13	2.00	0.43
1:W:30:LEU:HA	1:W:30:LEU:HD12	1.74	0.43
1:T:54:VAL:HG21	1:T:60:LEU:HD23	2.01	0.43
1:6:30:LEU:HD12	1:6:30:LEU:HA	1.76	0.43
1:U:19:ILE:HD12	1:U:31:LEU:HD22	2.01	0.43
1:X:5:LYS:HE3	1:Y:53:TRP:CD1	2.72	0.43
1:P:31:LEU:O	1:P:32:ASP:C	2.57	0.43
1:V:37:ALA:O	1:V:41:ILE:HB	2.18	0.43
1:U:7:ASN:OD1	1:U:9:ALA:HB3	2.24	0.43
1:L:153:ALA:C	1:L:154:ILE:HG12	2.45	0.43
1:4:142:LEU:HA	1:4:142:LEU:HD12	1.83	0.43
1:S:142:LEU:HD12	1:S:142:LEU:HA	1.86	0.43
1:9:56:GLY:O	1:9:59:GLU:HG2	2.18	0.43
1:P:9:ALA:HA	1:P:42:GLY:HA2	2.04	0.43
1:L:31:LEU:O	1:L:32:ASP:C	2.59	0.43
1:G:31:LEU:O	1:G:32:ASP:C	2.57	0.43
1:5:11:PRO:HB3	1:5:43:GLN:HB3	2.00	0.43
1:N:154:ILE:HD13	1:O:67:LEU:CD2	2.49	0.42
1:J:126:ILE:H	1:J:126:ILE:HG12	1.69	0.42
1:M:19:ILE:HD11	1:M:31:LEU:HB2	1.99	0.42
1:N:30:LEU:HA	1:N:30:LEU:HD12	1.73	0.42
1:R:142:LEU:HD12	1:R:142:LEU:HA	1.84	0.42
1:I:30:LEU:O	1:I:137:GLY:HA3	2.20	0.42
1:I:7:ASN:ND2	1:9:40:ARG:O	110.55	0.42
1:I:154:ILE:HD13	1:J:67:LEU:HD21	2.08	0.42
1:N:62:LEU:HD23	1:N:62:LEU:HA	1.81	0.42
1:8:54:VAL:HG11	1:8:63:ALA:HB2	2.00	0.42
1:E:31:LEU:O	1:E:32:ASP:C	2.58	0.42
1:T:108:SER:C	1:T:110:VAL:H	2.23	0.42
1:W:105:ALA:HB2	1:W:112:VAL:HG23	2.01	0.42
1:O:26:ILE:HD13	1:O:125:ALA:HB1	2.00	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:K:141:ALA:O	1:K:144:ALA:HB3	2.19	0.42
1:Z:62:LEU:HD13	1:4:151:LEU:HD11	2.01	0.42
1:D:60:LEU:HA	1:D:60:LEU:HD23	1.78	0.42
1:E:9:ALA:HA	1:E:42:GLY:HA2	2.01	0.42
1:R:31:LEU:O	1:R:34:ALA:N	2.57	0.42
1:S:62:LEU:HD23	1:S:62:LEU:HA	1.83	0.42
1:L:151:LEU:O	1:L:154:ILE:HG13	2.19	0.42
1:8:9:ALA:HA	1:8:42:GLY:HA2	2.00	0.42
1:D:54:VAL:HG11	1:D:63:ALA:HB2	2.07	0.42
1:7:153:ALA:O	1:7:154:ILE:HG12	2.19	0.42
1:5:118:THR:O	1:6:87:THR:HB	2.18	0.42
1:C:62:LEU:HA	1:C:62:LEU:HD23	1.79	0.42
1:E:36:ASP:OD2	1:E:40:ARG:NH1	2.52	0.42
1:U:147:MET:O	1:U:148:ILE:C	2.57	0.42
1:F:105:ALA:HB2	1:F:112:VAL:HG23	2.09	0.42
1:M:131:THR:HG23	1:X:25:PHE:HB3	80.42	0.42
1:O:18:THR:HG21	1:O:64:THR:HG22	2.01	0.42
1:K:18:THR:OG1	1:K:64:THR:HG22	2.19	0.42
1:H:30:LEU:O	1:H:137:GLY:HA3	2.23	0.42
1:W:7:ASN:OD1	1:W:9:ALA:HB3	2.18	0.42
1:F:41:ILE:HA	1:Q:7:ASN:ND2	2.33	0.42
1:L:54:VAL:HG11	1:L:63:ALA:HB2	2.06	0.42
1:V:50:THR:CG2	1:V:73:TYR:HE1	2.37	0.42
1:T:19:ILE:HG12	1:T:79:LEU:CB	2.51	0.42
1:P:151:LEU:O	1:P:154:ILE:HG13	2.19	0.42
1:I:7:ASN:OD1	1:I:9:ALA:HB3	2.21	0.42
1:F:7:ASN:ND2	1:F:8:VAL:H	2.24	0.42
1:8:55:PRO:HG2	1:8:59:GLU:HG3	2.00	0.42
1:2:34:ALA:HA	1:2:138:ALA:HA	2.01	0.42
1:L:84:ARG:HH21	1:L:90:PHE:HE1	1.80	0.42
1:J:27:ASN:HA	1:J:30:LEU:HB2	2.01	0.42
1:J:30:LEU:HD12	1:J:30:LEU:HA	1.75	0.42
1:R:126:ILE:HG12	1:R:126:ILE:H	1.71	0.42
1:X:83:ILE:HD13	1:X:122:ILE:HA	2.10	0.42
1:4:30:LEU:O	1:4:137:GLY:HA3	2.19	0.42
1:D:131:THR:HA	1:F:23:ASN:OD1	2.31	0.42
1:R:22:PHE:HD1	1:R:23:ASN:N	2.17	0.42
1:J:62:LEU:HA	1:J:62:LEU:HD23	1.85	0.42
1:C:31:LEU:O	1:C:32:ASP:C	2.62	0.42
1:2:50:THR:CG2	1:2:73:TYR:HE1	2.31	0.42
1:B:34:ALA:HA	1:B:138:ALA:HA	2.05	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:4:54:VAL:HB	1:4:55:PRO:HD2	2.01	0.42
1:H:89:HIS:HB3	2:H:157:SO4:O2	2.19	0.42
1:X:18:THR:HG21	1:X:64:THR:HG22	2.03	0.42
1:V:30:LEU:HA	1:V:30:LEU:HD12	1.78	0.42
1:9:30:LEU:HA	1:9:30:LEU:HD12	1.74	0.42
1:M:142:LEU:HA	1:M:142:LEU:HD12	1.86	0.42
1:K:37:ALA:O	1:K:41:ILE:HB	2.25	0.42
1:N:37:ALA:O	1:N:41:ILE:HB	2.24	0.42
1:I:41:ILE:HD13	1:O:7:ASN:ND2	2.34	0.42
1:R:19:ILE:HD11	1:R:31:LEU:HB2	2.01	0.42
1:R:31:LEU:O	1:R:32:ASP:C	2.58	0.42
1:R:9:ALA:HA	1:R:42:GLY:HA2	2.02	0.42
1:P:151:LEU:HD23	1:P:151:LEU:HA	1.90	0.42
1:Q:62:LEU:HD23	1:Q:62:LEU:HA	1.81	0.42
1:Z:9:ALA:HA	1:Z:42:GLY:HA2	2.01	0.42
1:4:54:VAL:HG11	1:4:63:ALA:HB2	2.02	0.42
1:7:26:ILE:HD13	1:7:125:ALA:HB1	2.00	0.42
1:D:18:THR:HG21	1:D:64:THR:HG22	2.00	0.42
1:N:153:ALA:C	1:N:154:ILE:HG12	2.42	0.42
1:2:18:THR:HG21	1:2:64:THR:HG22	2.02	0.42
1:9:111:PRO:HG3	1:9:151:LEU:HD12	2.01	0.42
1:U:22:PHE:HD1	1:U:23:ASN:N	2.17	0.42
1:A:9:ALA:HA	1:A:42:GLY:HA2	2.07	0.42
1:O:151:LEU:O	1:O:154:ILE:HG13	2.21	0.42
1:P:154:ILE:HD13	1:Q:67:LEU:CD2	2.55	0.42
1:I:62:LEU:HA	1:I:62:LEU:HD23	1.79	0.42
1:C:19:ILE:HD12	1:C:31:LEU:HD22	2.15	0.42
1:Z:20:ALA:HA	1:Z:54:VAL:O	2.19	0.42
1:A:54:VAL:HG11	1:A:63:ALA:HB2	2.02	0.42
1:A:55:PRO:HB2	1:E:143:THR:HG22	2.03	0.42
1:V:31:LEU:O	1:V:32:ASP:C	2.59	0.42
1:Z:67:LEU:HD21	1:4:154:ILE:HD13	2.00	0.42
1:A:119:THR:HG22	1:B:87:THR:HG21	2.00	0.42
1:8:118:THR:O	1:9:87:THR:HB	2.19	0.42
1:4:36:ASP:OD2	1:4:40:ARG:NH1	2.52	0.42
1:U:60:LEU:HD23	1:U:60:LEU:HA	1.89	0.42
1:4:27:ASN:HA	1:4:30:LEU:HB2	2.01	0.42
1:O:142:LEU:HD12	1:O:142:LEU:HA	1.86	0.42
1:R:153:ALA:C	1:R:154:ILE:HG12	2.47	0.42
1:W:62:LEU:HA	1:W:62:LEU:HD23	1.82	0.42
1:8:7:ASN:OD1	1:8:9:ALA:HB3	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:K:62:LEU:HD13	1:O:151:LEU:HD11	2.05	0.42
1:J:31:LEU:O	1:J:34:ALA:N	2.60	0.42
1:A:31:LEU:O	1:A:32:ASP:C	2.57	0.42
1:C:20:ALA:HA	1:C:54:VAL:O	2.19	0.42
1:E:34:ALA:HA	1:E:138:ALA:HA	2.01	0.42
1:F:85:GLY:HA3	2:F:157:SO4:O3	2.29	0.42
1:R:62:LEU:HD23	1:R:62:LEU:HA	1.84	0.42
1:G:25:PHE:CE1	1:G:126:ILE:HG23	5.20	0.42
1:T:18:THR:HG21	1:T:64:THR:HG22	2.01	0.42
1:M:123:GLU:OE2	1:8:122:ILE:HD13	53.91	0.42
1:C:2:ASN:HD21	1:D:50:THR:CG2	2.45	0.42
1:P:54:VAL:HG11	1:P:63:ALA:HB2	2.02	0.42
1:5:87:THR:HG21	1:9:119:THR:HG22	2.00	0.42
1:M:7:ASN:ND2	1:7:41:ILE:HA	100.30	0.42
1:I:19:ILE:HD12	1:I:31:LEU:HD22	2.02	0.42
1:P:153:ALA:C	1:P:154:ILE:HG12	2.43	0.42
1:L:1:MET:HB2	1:M:46:ASP:O	2.19	0.42
1:1:1:MET:HA	1:2:46:ASP:O	2.19	0.42
1:6:111:PRO:HG3	1:6:151:LEU:HD12	2.01	0.42
1:N:141:ALA:O	1:N:144:ALA:HB3	2.28	0.42
1:F:60:LEU:HA	1:F:60:LEU:HD23	1.80	0.42
1:V:60:LEU:HA	1:V:60:LEU:HD23	1.86	0.42
1:7:5:LYS:CE	1:8:21:ARG:HH22	2.25	0.42
1:Q:30:LEU:O	1:Q:137:GLY:HA3	2.19	0.42
1:2:22:PHE:HD1	1:2:23:ASN:N	2.18	0.42
1:Y:30:LEU:HD12	1:Y:30:LEU:HA	1.75	0.42
1:P:19:ILE:HD12	1:P:31:LEU:HD22	2.04	0.42
1:H:62:LEU:HD23	1:H:62:LEU:HA	1.78	0.42
1:N:142:LEU:HD12	1:N:142:LEU:HA	1.85	0.42
1:E:7:ASN:ND2	1:N:41:ILE:HA	2.35	0.42
1:M:67:LEU:O	1:M:70:SER:HB3	2.20	0.42
1:H:151:LEU:O	1:H:154:ILE:HG13	2.20	0.42
1:J:31:LEU:O	1:J:32:ASP:C	2.60	0.42
1:U:67:LEU:CD2	1:Y:154:ILE:HD13	2.50	0.42
1:7:31:LEU:O	1:7:32:ASP:C	2.58	0.42
1:7:84:ARG:HH21	1:7:90:PHE:HE1	1.68	0.42
1:Z:34:ALA:HA	1:Z:138:ALA:HA	2.01	0.42
1:B:64:THR:HB	1:B:76:VAL:HG11	2.01	0.42
1:1:122:ILE:CD1	1:1:122:ILE:H	2.15	0.42
1:F:151:LEU:HD23	1:F:151:LEU:HA	1.92	0.42
1:O:22:PHE:HD1	1:O:23:ASN:N	2.18	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:P:19:ILE:HD11	1:P:31:LEU:HB2	2.02	0.42
1:A:41:ILE:HD12	1:L:8:VAL:HG22	73.94	0.42
1:G:7:ASN:OD1	1:G:9:ALA:HB3	2.20	0.42
1:H:142:LEU:HD12	1:H:142:LEU:HA	1.82	0.42
1:W:37:ALA:O	1:W:41:ILE:HB	2.19	0.42
1:H:9:ALA:HA	1:H:42:GLY:HA2	2.02	0.42
1:Z:7:ASN:OD1	1:Z:9:ALA:HB3	2.19	0.42
1:I:7:ASN:ND2	1:5:41:ILE:HA	2.34	0.42
1:L:34:ALA:HA	1:L:138:ALA:HA	2.01	0.42
1:I:60:LEU:HA	1:I:60:LEU:HD23	1.89	0.41
1:9:18:THR:HG21	1:9:64:THR:HG22	2.01	0.41
1:N:25:PHE:CE1	1:W:126:ILE:HG23	60.48	0.41
1:R:5:LYS:NZ	1:S:21:ARG:NH2	2.72	0.41
1:P:83:ILE:HD13	1:P:122:ILE:HA	2.02	0.41
1:M:30:LEU:HA	1:M:30:LEU:HD12	1.77	0.41
1:Q:27:ASN:HA	1:Q:30:LEU:HB2	2.04	0.41
1:R:40:ARG:CZ	1:X:21:ARG:CZ	92.45	0.41
1:L:9:ALA:HA	1:L:42:GLY:HA2	2.02	0.41
1:Z:36:ASP:OD2	1:Z:40:ARG:NH1	2.53	0.41
1:O:62:LEU:HD23	1:O:62:LEU:HA	1.79	0.41
1:8:22:PHE:HD1	1:8:23:ASN:N	2.17	0.41
1:B:27:ASN:HA	1:B:30:LEU:HB2	2.01	0.41
1:X:5:LYS:CE	1:Y:21:ARG:HH22	2.30	0.41
1:Y:30:LEU:O	1:Y:137:GLY:HA3	2.20	0.41
1:Y:19:ILE:HD12	1:Y:31:LEU:HD22	2.02	0.41
1:D:130:GLY:O	1:F:23:ASN:HA	2.30	0.41
1:B:37:ALA:O	1:B:41:ILE:HB	2.20	0.41
1:A:142:LEU:HA	1:A:142:LEU:HD12	1.88	0.41
1:I:31:LEU:O	1:I:32:ASP:C	2.58	0.41
1:J:19:ILE:HD12	1:J:31:LEU:HD22	2.19	0.41
1:8:2:ASN:HD21	1:9:50:THR:HG22	1.85	0.41
1:4:111:PRO:HG3	1:4:151:LEU:HD12	2.01	0.41
1:J:26:ILE:HD13	1:J:125:ALA:HB1	2.09	0.41
1:X:7:ASN:ND2	1:4:41:ILE:HA	2.35	0.41
1:T:26:ILE:HD13	1:T:125:ALA:HB1	2.13	0.41
1:Y:105:ALA:HB2	1:Y:112:VAL:HG23	2.05	0.41
1:J:18:THR:HG21	1:J:64:THR:HG22	2.04	0.41
1:E:67:LEU:O	1:E:70:SER:HB3	2.21	0.41
1:A:18:THR:HG21	1:A:64:THR:HG22	2.02	0.41
1:N:153:ALA:C	1:N:154:ILE:CG1	2.91	0.41
1:O:18:THR:OG1	1:O:64:THR:HG22	2.31	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:N:31:LEU:O	1:N:34:ALA:N	2.62	0.41
1:E:7:ASN:ND2	1:E:8:VAL:H	2.18	0.41
1:D:142:LEU:HD12	1:D:142:LEU:HA	1.91	0.41
1:W:16:ALA:CB	1:W:67:LEU:HD13	2.49	0.41
1:V:31:LEU:O	1:V:35:VAL:HG23	2.31	0.41
1:2:19:ILE:HG21	1:2:53:TRP:CZ3	2.55	0.41
1:E:141:ALA:O	1:E:144:ALA:HB3	2.32	0.41
1:F:31:LEU:O	1:F:34:ALA:N	2.60	0.41
1:W:5:LYS:CE	1:X:21:ARG:HH22	2.28	0.41
1:X:31:LEU:O	1:X:34:ALA:N	2.62	0.41
1:O:31:LEU:O	1:O:34:ALA:N	2.56	0.41
1:W:2:ASN:HD21	1:X:50:THR:CG2	2.45	0.41
1:2:27:ASN:HA	1:2:30:LEU:HB2	2.02	0.41
1:P:142:LEU:HA	1:P:142:LEU:HD12	1.86	0.41
1:C:50:THR:CG2	1:C:73:TYR:HE1	2.34	0.41
1:Q:111:PRO:HG3	1:Q:151:LEU:HD12	2.01	0.41
1:E:153:ALA:C	1:E:154:ILE:HG12	2.51	0.41
1:C:153:ALA:O	1:C:154:ILE:HG12	2.20	0.41
1:6:9:ALA:HA	1:6:42:GLY:HA2	2.01	0.41
1:B:62:LEU:HA	1:B:62:LEU:HD23	1.86	0.41
1:W:36:ASP:OD2	1:W:40:ARG:NH1	2.53	0.41
1:Q:31:LEU:O	1:Q:32:ASP:C	2.57	0.41
1:V:105:ALA:HB2	1:V:112:VAL:HG23	2.10	0.41
1:J:105:ALA:HB2	1:J:112:VAL:HG23	2.15	0.41
1:K:52:VAL:HG22	1:O:4:ILE:HD12	2.02	0.41
1:I:34:ALA:HA	1:I:138:ALA:HA	2.01	0.41
1:W:126:ILE:HG12	1:W:126:ILE:H	1.71	0.41
1:W:83:ILE:HD13	1:W:122:ILE:HA	2.02	0.41
1:S:126:ILE:HG12	1:S:126:ILE:H	1.72	0.41
1:F:19:ILE:HD12	1:F:31:LEU:HD22	2.04	0.41
1:F:30:LEU:HD12	1:F:30:LEU:HA	1.77	0.41
1:P:53:TRP:CD1	1:T:5:LYS:HE3	2.56	0.41
1:O:31:LEU:O	1:O:35:VAL:HG23	2.19	0.41
1:B:7:ASN:ND2	1:B:8:VAL:H	2.25	0.41
1:T:37:ALA:O	1:T:41:ILE:HB	2.25	0.41
1:M:118:THR:O	1:N:87:THR:HB	2.45	0.41
1:9:36:ASP:OD2	1:9:40:ARG:NH1	2.53	0.41
1:H:7:ASN:ND2	1:H:8:VAL:H	2.19	0.41
1:5:34:ALA:HA	1:5:138:ALA:HA	2.02	0.41
1:5:7:ASN:OD1	1:5:9:ALA:HB3	2.20	0.41
1:Y:62:LEU:HD23	1:Y:62:LEU:HA	1.82	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:50:THR:CG2	1:H:73:TYR:HE1	2.43	0.41
1:P:50:THR:CG2	1:T:2:ASN:HD21	2.31	0.41
1:O:111:PRO:HG3	1:O:151:LEU:HD12	2.07	0.41
1:P:62:LEU:HD13	1:T:151:LEU:HD11	2.03	0.41
1:S:36:ASP:OD2	1:S:40:ARG:NH1	2.53	0.41
1:K:34:ALA:HA	1:K:138:ALA:HA	2.05	0.41
1:Q:105:ALA:HB2	1:Q:112:VAL:HG23	2.03	0.41
1:3:26:ILE:HD13	1:3:125:ALA:HB1	2.03	0.41
1:R:60:LEU:HD23	1:R:60:LEU:HA	1.85	0.41
1:T:16:ALA:CB	1:T:67:LEU:HD13	2.55	0.41
1:3:27:ASN:HA	1:3:30:LEU:HB2	2.03	0.41
1:M:37:ALA:O	1:M:41:ILE:HB	2.21	0.41
1:H:5:LYS:CE	1:I:21:ARG:HH22	2.30	0.41
1:G:16:ALA:CB	1:G:67:LEU:HD13	2.53	0.41
1:P:27:ASN:HA	1:P:30:LEU:HB2	2.02	0.41
1:J:23:ASN:HA	1:9:130:GLY:O	83.32	0.41
1:T:7:ASN:ND2	1:T:8:VAL:H	2.28	0.41
1:H:20:ALA:HA	1:H:54:VAL:O	2.21	0.41
1:V:153:ALA:C	1:V:154:ILE:HG12	2.47	0.41
1:S:7:ASN:OD1	1:S:9:ALA:HB3	2.20	0.41
1:D:62:LEU:HD23	1:D:62:LEU:HA	1.87	0.41
1:Z:67:LEU:CD2	1:4:154:ILE:HD13	2.50	0.41
1:C:54:VAL:HG11	1:C:63:ALA:HB2	2.02	0.41
1:W:31:LEU:O	1:W:34:ALA:N	2.57	0.41
1:E:31:LEU:O	1:E:34:ALA:N	2.53	0.41
1:8:34:ALA:HA	1:8:138:ALA:HA	2.02	0.41
1:N:151:LEU:O	1:N:154:ILE:HG13	2.21	0.41
1:2:16:ALA:O	1:2:76:VAL:HA	2.21	0.41
1:J:83:ILE:HD13	1:J:122:ILE:HA	2.03	0.41
1:E:21:ARG:O	1:E:24:GLN:HB3	2.21	0.41
1:C:9:ALA:HA	1:C:42:GLY:HA2	2.04	0.41
1:T:153:ALA:C	1:T:154:ILE:CG1	2.92	0.41
1:R:19:ILE:CG2	1:R:53:TRP:CZ3	3.05	0.41
1:O:151:LEU:HA	1:O:151:LEU:HD23	2.03	0.41
1:S:7:ASN:ND2	1:Y:40:ARG:O	2.54	0.41
1:2:154:ILE:HD13	1:3:67:LEU:CD2	2.51	0.41
1:R:20:ALA:HA	1:R:54:VAL:O	2.21	0.41
1:B:31:LEU:O	1:B:32:ASP:C	2.59	0.41
1:I:153:ALA:C	1:I:154:ILE:HG12	2.40	0.41
1:R:141:ALA:O	1:R:144:ALA:HB3	2.20	0.41
1:G:23:ASN:HA	1:Q:131:THR:HA	2.03	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:K:126:ILE:HG12	1:K:126:ILE:H	1.67	0.41
1:V:5:LYS:CE	1:W:21:ARG:NH2	2.80	0.41
1:P:85:GLY:CA	2:P:157:SO4:O4	2.63	0.41
1:F:154:ILE:HD13	1:G:67:LEU:HD21	2.12	0.41
1:9:22:PHE:HD1	1:9:23:ASN:N	2.18	0.41
1:N:22:PHE:HE1	1:7:131:THR:O	75.05	0.41
1:K:131:THR:HA	1:M:23:ASN:HA	46.59	0.41
1:K:130:GLY:O	1:M:23:ASN:HA	47.59	0.41
1:Y:145:LEU:O	1:Y:146:GLU:C	2.59	0.41
1:P:2:ASN:HD21	1:Q:50:THR:CG2	2.34	0.41
1:2:7:ASN:OD1	1:2:9:ALA:HB3	2.20	0.41
1:G:8:VAL:HG22	1:P:41:ILE:HD12	45.42	0.41
1:V:154:ILE:HD13	1:W:67:LEU:CD2	2.61	0.41
1:Q:151:LEU:HD23	1:Q:151:LEU:HA	1.91	0.41
1:S:31:LEU:O	1:S:32:ASP:C	2.59	0.41
1:S:19:ILE:HD11	1:S:31:LEU:HB2	2.03	0.41
1:H:7:ASN:OD1	1:H:9:ALA:HB3	2.26	0.41
1:Q:40:ARG:O	1:Z:7:ASN:ND2	119.89	0.41
1:4:9:ALA:HA	1:4:42:GLY:HA2	2.01	0.41
1:U:67:LEU:O	1:U:70:SER:HB3	2.21	0.41
1:N:1:MET:HA	1:O:46:ASP:O	2.21	0.41
1:F:87:THR:HB	1:J:118:THR:O	2.21	0.41
1:J:84:ARG:HH21	1:J:90:PHE:HE1	1.79	0.41
1:R:26:ILE:HD13	1:R:125:ALA:HB1	2.02	0.41
1:B:141:ALA:O	1:B:144:ALA:HB3	2.27	0.41
1:5:31:LEU:O	1:5:32:ASP:C	2.59	0.41
1:E:60:LEU:HA	1:E:60:LEU:HD23	1.86	0.41
1:S:18:THR:HG21	1:S:64:THR:HG22	2.03	0.41
1:F:21:ARG:NH2	1:J:5:LYS:NZ	2.76	0.41
1:Q:142:LEU:HD12	1:Q:142:LEU:HA	1.94	0.41
1:I:142:LEU:HD12	1:I:142:LEU:HA	1.84	0.41
1:I:37:ALA:O	1:I:41:ILE:HB	2.25	0.41
1:L:151:LEU:HD23	1:L:151:LEU:HA	1.90	0.41
1:A:151:LEU:O	1:A:154:ILE:HG13	2.24	0.41
1:8:154:ILE:HD13	1:9:67:LEU:HD21	2.03	0.41
1:8:19:ILE:HD11	1:8:31:LEU:HB2	2.02	0.41
1:F:62:LEU:HD23	1:F:62:LEU:HA	1.84	0.41
1:F:46:ASP:O	1:J:1:MET:HB2	2.22	0.41
1:G:83:ILE:HD13	1:G:122:ILE:HA	2.03	0.40
1:E:27:ASN:HA	1:E:30:LEU:HB2	2.04	0.40
1:L:20:ALA:HA	1:L:54:VAL:O	2.21	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:V:142:LEU:HD12	1:V:142:LEU:HA	1.83	0.40
1:B:7:ASN:OD1	1:B:9:ALA:HB3	2.21	0.40
1:U:2:ASN:HD21	1:V:50:THR:CG2	2.41	0.40
1:O:153:ALA:C	1:O:154:ILE:HG12	2.42	0.40
1:7:9:ALA:HA	1:7:42:GLY:HA2	2.03	0.40
1:5:111:PRO:HG3	1:5:151:LEU:HD12	2.03	0.40
1:J:60:LEU:HD23	1:J:60:LEU:HA	1.86	0.40
1:C:60:LEU:HA	1:C:60:LEU:HD23	1.81	0.40
1:W:18:THR:HG21	1:W:64:THR:HG22	2.03	0.40
1:C:142:LEU:HA	1:C:142:LEU:HD12	1.84	0.40
1:M:143:THR:HG22	1:N:55:PRO:CB	2.50	0.40
1:Q:67:LEU:O	1:Q:70:SER:HB3	2.21	0.40
1:8:142:LEU:HD12	1:8:142:LEU:HA	1.82	0.40
1:C:19:ILE:CG2	1:C:53:TRP:CZ3	3.11	0.40
1:Y:7:ASN:ND2	1:Y:8:VAL:H	2.27	0.40
1:B:26:ILE:HD13	1:B:125:ALA:HB1	2.08	0.40
1:T:34:ALA:HA	1:T:138:ALA:HA	2.03	0.40
1:B:108:SER:C	1:B:110:VAL:H	2.31	0.40
1:U:34:ALA:HA	1:U:138:ALA:HA	2.02	0.40
1:I:18:THR:HG21	1:I:64:THR:HG22	2.03	0.40
1:D:111:PRO:HG3	1:D:151:LEU:HD12	2.02	0.40
1:E:16:ALA:O	1:E:76:VAL:HA	2.22	0.40
1:G:22:PHE:HE1	1:G:131:THR:O	22.41	0.40
1:C:5:LYS:HE3	1:D:53:TRP:CD1	2.69	0.40
1:V:18:THR:HG21	1:V:64:THR:HG22	2.07	0.40
1:N:83:ILE:HD13	1:N:122:ILE:HA	2.03	0.40
1:H:130:GLY:O	1:3:23:ASN:HA	71.31	0.40
1:K:142:LEU:HA	1:K:142:LEU:HD12	1.85	0.40
1:U:153:ALA:C	1:U:154:ILE:HG12	2.42	0.40
1:S:16:ALA:CB	1:S:67:LEU:HD13	2.50	0.40
1:T:7:ASN:OD1	1:T:9:ALA:HB3	2.21	0.40
1:C:36:ASP:OD2	1:C:40:ARG:NH1	2.53	0.40
1:S:50:THR:CG2	1:S:73:TYR:HE1	2.35	0.40
1:C:19:ILE:HD11	1:C:31:LEU:HB2	2.03	0.40
1:M:151:LEU:HD11	1:N:62:LEU:HD13	2.15	0.40
1:1:54:VAL:HG11	1:1:63:ALA:HB2	2.02	0.40
1:4:151:LEU:HD23	1:4:151:LEU:HA	1.90	0.40
1:T:31:LEU:O	1:T:34:ALA:N	2.57	0.40
1:C:118:THR:O	1:D:87:THR:HB	2.28	0.40
1:5:26:ILE:HD13	1:5:125:ALA:HB1	2.03	0.40
1:B:105:ALA:HB2	1:B:112:VAL:HG23	2.13	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:1:18:THR:HG21	1:1:64:THR:HG22	2.04	0.40
1:1:27:ASN:HA	1:1:30:LEU:HB2	2.04	0.40
1:F:153:ALA:C	1:F:154:ILE:CG1	2.90	0.40
1:T:22:PHE:HD1	1:T:23:ASN:H	1.78	0.40
1:Z:22:PHE:HD1	1:Z:23:ASN:N	2.19	0.40
1:F:142:LEU:HA	1:F:142:LEU:HD12	1.91	0.40
1:A:7:ASN:ND2	1:A:8:VAL:H	2.19	0.40
1:L:37:ALA:O	1:L:41:ILE:HB	2.21	0.40
1:K:119:THR:HA	1:L:87:THR:HB	2.20	0.40
1:1:142:LEU:HD12	1:1:142:LEU:HA	1.86	0.40
1:R:16:ALA:CB	1:R:67:LEU:HD13	2.50	0.40
1:7:151:LEU:O	1:7:154:ILE:HG13	2.21	0.40
1:1:151:LEU:O	1:1:154:ILE:HG13	2.21	0.40
1:P:36:ASP:OD2	1:P:40:ARG:NH1	2.55	0.40
1:L:26:ILE:HD13	1:L:125:ALA:HB1	2.03	0.40
1:A:26:ILE:HD13	1:A:125:ALA:HB1	2.05	0.40
1:Q:34:ALA:HA	1:Q:138:ALA:HA	2.06	0.40
1:M:131:THR:HA	1:8:23:ASN:OD1	70.44	0.40
1:D:151:LEU:HD23	1:D:151:LEU:HA	1.92	0.40
1:D:151:LEU:O	1:D:154:ILE:HG13	2.21	0.40
1:O:16:ALA:O	1:O:76:VAL:HA	2.26	0.40
1:3:30:LEU:O	1:3:137:GLY:HA3	2.21	0.40
1:X:19:ILE:HD11	1:X:31:LEU:HB2	2.08	0.40
1:X:36:ASP:OD2	1:X:40:ARG:NH1	2.53	0.40
1:W:151:LEU:O	1:W:154:ILE:HG13	2.24	0.40
1:J:7:ASN:OD1	1:J:9:ALA:HB3	2.21	0.40
1:B:61:PRO:O	1:B:62:LEU:C	2.60	0.40
1:M:141:ALA:O	1:M:144:ALA:HB3	2.24	0.40
1:C:34:ALA:HA	1:C:138:ALA:HA	2.04	0.40
1:F:108:SER:C	1:F:110:VAL:H	2.29	0.40
1:Z:46:ASP:O	1:4:1:MET:HB2	2.22	0.40

There are no symmetry-related clashes.

## 5.3 Torsion angles ⓘ

### 5.3.1 Protein backbone ⓘ

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	1	152/156 (97%)	131 (86%)	19 (12%)	2 (1%)	15	61
1	2	152/156 (97%)	132 (87%)	18 (12%)	2 (1%)	15	61
1	3	152/156 (97%)	133 (88%)	17 (11%)	2 (1%)	15	61
1	4	152/156 (97%)	132 (87%)	18 (12%)	2 (1%)	15	61
1	5	152/156 (97%)	133 (88%)	17 (11%)	2 (1%)	15	61
1	6	152/156 (97%)	131 (86%)	19 (12%)	2 (1%)	15	61
1	7	152/156 (97%)	132 (87%)	18 (12%)	2 (1%)	15	61
1	8	152/156 (97%)	133 (88%)	17 (11%)	2 (1%)	15	61
1	9	152/156 (97%)	132 (87%)	18 (12%)	2 (1%)	15	61
1	A	152/156 (97%)	131 (86%)	19 (12%)	2 (1%)	15	61
1	B	152/156 (97%)	131 (86%)	19 (12%)	2 (1%)	15	61
1	C	152/156 (97%)	130 (86%)	19 (12%)	3 (2%)	9	53
1	D	152/156 (97%)	133 (88%)	17 (11%)	2 (1%)	15	61
1	E	152/156 (97%)	131 (86%)	19 (12%)	2 (1%)	15	61
1	F	152/156 (97%)	133 (88%)	17 (11%)	2 (1%)	15	61
1	G	152/156 (97%)	131 (86%)	19 (12%)	2 (1%)	15	61
1	H	152/156 (97%)	134 (88%)	16 (10%)	2 (1%)	15	61
1	I	152/156 (97%)	131 (86%)	19 (12%)	2 (1%)	15	61
1	J	152/156 (97%)	132 (87%)	18 (12%)	2 (1%)	15	61
1	K	152/156 (97%)	130 (86%)	20 (13%)	2 (1%)	15	61
1	L	152/156 (97%)	131 (86%)	19 (12%)	2 (1%)	15	61
1	M	152/156 (97%)	131 (86%)	19 (12%)	2 (1%)	15	61
1	N	152/156 (97%)	131 (86%)	19 (12%)	2 (1%)	15	61
1	O	152/156 (97%)	131 (86%)	19 (12%)	2 (1%)	15	61
1	P	152/156 (97%)	131 (86%)	19 (12%)	2 (1%)	15	61
1	Q	152/156 (97%)	130 (86%)	20 (13%)	2 (1%)	15	61
1	R	152/156 (97%)	132 (87%)	18 (12%)	2 (1%)	15	61
1	S	152/156 (97%)	132 (87%)	18 (12%)	2 (1%)	15	61
1	T	152/156 (97%)	131 (86%)	19 (12%)	2 (1%)	15	61
1	U	152/156 (97%)	133 (88%)	17 (11%)	2 (1%)	15	61

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	V	152/156 (97%)	131 (86%)	19 (12%)	2 (1%)	15	61
1	W	152/156 (97%)	131 (86%)	19 (12%)	2 (1%)	15	61
1	X	152/156 (97%)	131 (86%)	19 (12%)	2 (1%)	15	61
1	Y	152/156 (97%)	133 (88%)	17 (11%)	2 (1%)	15	61
1	Z	152/156 (97%)	133 (88%)	17 (11%)	2 (1%)	15	61
1	a	152/156 (97%)	132 (87%)	18 (12%)	2 (1%)	15	61
1	b	152/156 (97%)	133 (88%)	17 (11%)	2 (1%)	15	61
1	c	152/156 (97%)	130 (86%)	20 (13%)	2 (1%)	15	61
1	d	152/156 (97%)	132 (87%)	18 (12%)	2 (1%)	15	61
1	e	152/156 (97%)	133 (88%)	16 (10%)	3 (2%)	9	53
1	f	152/156 (97%)	131 (86%)	19 (12%)	2 (1%)	15	61
1	g	152/156 (97%)	129 (85%)	21 (14%)	2 (1%)	15	61
1	h	152/156 (97%)	131 (86%)	19 (12%)	2 (1%)	15	61
1	i	152/156 (97%)	131 (86%)	19 (12%)	2 (1%)	15	61
1	j	152/156 (97%)	133 (88%)	17 (11%)	2 (1%)	15	61
1	k	152/156 (97%)	132 (87%)	18 (12%)	2 (1%)	15	61
1	l	152/156 (97%)	133 (88%)	17 (11%)	2 (1%)	15	61
1	m	152/156 (97%)	133 (88%)	17 (11%)	2 (1%)	15	61
1	n	152/156 (97%)	133 (88%)	17 (11%)	2 (1%)	15	61
1	o	152/156 (97%)	132 (87%)	18 (12%)	2 (1%)	15	61
1	p	152/156 (97%)	131 (86%)	19 (12%)	2 (1%)	15	61
1	q	152/156 (97%)	131 (86%)	19 (12%)	2 (1%)	15	61
1	r	152/156 (97%)	129 (85%)	21 (14%)	2 (1%)	15	61
1	s	152/156 (97%)	131 (86%)	19 (12%)	2 (1%)	15	61
1	t	152/156 (97%)	134 (88%)	16 (10%)	2 (1%)	15	61
1	u	152/156 (97%)	130 (86%)	20 (13%)	2 (1%)	15	61
1	v	152/156 (97%)	132 (87%)	18 (12%)	2 (1%)	15	61
1	w	152/156 (97%)	131 (86%)	19 (12%)	2 (1%)	15	61
1	x	152/156 (97%)	132 (87%)	18 (12%)	2 (1%)	15	61
1	y	152/156 (97%)	130 (86%)	19 (12%)	3 (2%)	9	53
All	All	9120/9360 (97%)	7897 (87%)	1100 (12%)	123 (1%)	15	61

All (123) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	I	61	PRO
1	J	61	PRO
1	Q	61	PRO
1	A	55	PRO
1	A	61	PRO
1	D	55	PRO
1	D	61	PRO
1	E	61	PRO
1	I	55	PRO
1	K	61	PRO
1	L	61	PRO
1	M	55	PRO
1	M	61	PRO
1	N	61	PRO
1	O	61	PRO
1	P	61	PRO
1	Q	55	PRO
1	S	61	PRO
1	T	61	PRO
1	V	61	PRO
1	W	55	PRO
1	Y	55	PRO
1	1	61	PRO
1	3	61	PRO
1	4	55	PRO
1	9	55	PRO
1	9	61	PRO
1	a	61	PRO
1	c	61	PRO
1	e	152	LYS
1	f	152	LYS
1	g	55	PRO
1	g	61	PRO
1	h	61	PRO
1	j	61	PRO
1	l	55	PRO
1	o	61	PRO
1	q	61	PRO
1	r	61	PRO
1	t	55	PRO
1	u	61	PRO
1	v	61	PRO

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Mol	Chain	Res	Type
1	B	55	PRO
1	B	61	PRO
1	C	55	PRO
1	C	61	PRO
1	C	152	LYS
1	F	55	PRO
1	F	61	PRO
1	G	61	PRO
1	H	55	PRO
1	L	55	PRO
1	P	55	PRO
1	S	55	PRO
1	U	55	PRO
1	V	55	PRO
1	W	61	PRO
1	X	61	PRO
1	Y	61	PRO
1	Z	55	PRO
1	1	55	PRO
1	2	55	PRO
1	2	61	PRO
1	4	61	PRO
1	5	55	PRO
1	6	61	PRO
1	7	55	PRO
1	a	55	PRO
1	d	55	PRO
1	d	61	PRO
1	i	55	PRO
1	i	61	PRO
1	j	55	PRO
1	l	61	PRO
1	m	55	PRO
1	o	55	PRO
1	p	55	PRO
1	p	61	PRO
1	q	55	PRO
1	s	55	PRO
1	s	61	PRO
1	v	55	PRO
1	w	61	PRO
1	x	55	PRO

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Mol	Chain	Res	Type
1	x	61	PRO
1	y	61	PRO
1	y	152	LYS
1	E	55	PRO
1	H	61	PRO
1	J	55	PRO
1	K	55	PRO
1	N	55	PRO
1	R	61	PRO
1	U	61	PRO
1	Z	61	PRO
1	b	55	PRO
1	b	61	PRO
1	e	61	PRO
1	f	55	PRO
1	n	55	PRO
1	n	61	PRO
1	r	55	PRO
1	t	61	PRO
1	u	55	PRO
1	G	55	PRO
1	O	55	PRO
1	T	55	PRO
1	X	55	PRO
1	3	55	PRO
1	6	55	PRO
1	7	61	PRO
1	c	55	PRO
1	h	55	PRO
1	k	61	PRO
1	w	55	PRO
1	y	55	PRO
1	R	55	PRO
1	5	61	PRO
1	8	61	PRO
1	e	55	PRO
1	k	55	PRO
1	m	61	PRO
1	8	55	PRO

### 5.3.2 Protein sidechains ⓘ

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	1	110/112 (98%)	95 (86%)	15 (14%)	5	28
1	2	109/112 (97%)	94 (86%)	15 (14%)	4	28
1	3	109/112 (97%)	94 (86%)	15 (14%)	4	28
1	4	109/112 (97%)	93 (85%)	16 (15%)	4	25
1	5	110/112 (98%)	95 (86%)	15 (14%)	5	28
1	6	110/112 (98%)	96 (87%)	14 (13%)	5	30
1	7	109/112 (97%)	95 (87%)	14 (13%)	5	30
1	8	109/112 (97%)	94 (86%)	15 (14%)	4	28
1	9	110/112 (98%)	94 (86%)	16 (14%)	4	26
1	A	109/112 (97%)	95 (87%)	14 (13%)	5	30
1	B	109/112 (97%)	93 (85%)	16 (15%)	4	25
1	C	109/112 (97%)	94 (86%)	15 (14%)	4	28
1	D	110/112 (98%)	95 (86%)	15 (14%)	5	28
1	E	109/112 (97%)	94 (86%)	15 (14%)	4	28
1	F	109/112 (97%)	95 (87%)	14 (13%)	5	30
1	G	109/112 (97%)	95 (87%)	14 (13%)	5	30
1	H	109/112 (97%)	93 (85%)	16 (15%)	4	25
1	I	109/112 (97%)	96 (88%)	13 (12%)	6	33
1	J	109/112 (97%)	93 (85%)	16 (15%)	4	25
1	K	109/112 (97%)	94 (86%)	15 (14%)	4	28
1	L	109/112 (97%)	94 (86%)	15 (14%)	4	28
1	M	109/112 (97%)	93 (85%)	16 (15%)	4	25
1	N	109/112 (97%)	94 (86%)	15 (14%)	4	28
1	O	109/112 (97%)	95 (87%)	14 (13%)	5	30
1	P	108/112 (96%)	93 (86%)	15 (14%)	4	28
1	Q	109/112 (97%)	93 (85%)	16 (15%)	4	25

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	R	109/112 (97%)	95 (87%)	14 (13%)	5	30
1	S	109/112 (97%)	96 (88%)	13 (12%)	6	33
1	T	109/112 (97%)	94 (86%)	15 (14%)	4	28
1	U	109/112 (97%)	94 (86%)	15 (14%)	4	28
1	V	109/112 (97%)	93 (85%)	16 (15%)	4	25
1	W	109/112 (97%)	95 (87%)	14 (13%)	5	30
1	X	109/112 (97%)	93 (85%)	16 (15%)	4	25
1	Y	108/112 (96%)	94 (87%)	14 (13%)	5	30
1	Z	109/112 (97%)	94 (86%)	15 (14%)	4	28
1	a	109/112 (97%)	95 (87%)	14 (13%)	5	30
1	b	110/112 (98%)	97 (88%)	13 (12%)	6	34
1	c	109/112 (97%)	94 (86%)	15 (14%)	4	28
1	d	109/112 (97%)	93 (85%)	16 (15%)	4	25
1	e	109/112 (97%)	93 (85%)	16 (15%)	4	25
1	f	110/112 (98%)	95 (86%)	15 (14%)	5	28
1	g	110/112 (98%)	94 (86%)	16 (14%)	4	26
1	h	109/112 (97%)	96 (88%)	13 (12%)	6	33
1	i	109/112 (97%)	94 (86%)	15 (14%)	4	28
1	j	110/112 (98%)	95 (86%)	15 (14%)	5	28
1	k	109/112 (97%)	93 (85%)	16 (15%)	4	25
1	l	110/112 (98%)	95 (86%)	15 (14%)	5	28
1	m	109/112 (97%)	94 (86%)	15 (14%)	4	28
1	n	109/112 (97%)	94 (86%)	15 (14%)	4	28
1	o	108/112 (96%)	93 (86%)	15 (14%)	4	28
1	p	109/112 (97%)	95 (87%)	14 (13%)	5	30
1	q	109/112 (97%)	95 (87%)	14 (13%)	5	30
1	r	109/112 (97%)	95 (87%)	14 (13%)	5	30
1	s	109/112 (97%)	94 (86%)	15 (14%)	4	28
1	t	109/112 (97%)	94 (86%)	15 (14%)	4	28
1	u	108/112 (96%)	93 (86%)	15 (14%)	4	28
1	v	109/112 (97%)	95 (87%)	14 (13%)	5	30

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	w	109/112 (97%)	94 (86%)	15 (14%)	4	28
1	x	108/112 (96%)	94 (87%)	14 (13%)	5	30
1	y	110/112 (98%)	96 (87%)	14 (13%)	5	30
All	All	6546/6720 (97%)	5657 (86%)	889 (14%)	5	28

All (889) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	2	ASN
1	A	19	ILE
1	A	22	PHE
1	A	24	GLN
1	A	43	GLN
1	A	45	LYS
1	A	50	THR
1	A	59	GLU
1	A	62	LEU
1	A	79	LEU
1	A	118	THR
1	A	136	LYS
1	A	142	LEU
1	A	154	ILE
1	B	2	ASN
1	B	19	ILE
1	B	22	PHE
1	B	24	GLN
1	B	30	LEU
1	B	43	GLN
1	B	45	LYS
1	B	50	THR
1	B	62	LEU
1	B	64	THR
1	B	79	LEU
1	B	116	VAL
1	B	118	THR
1	B	136	LYS
1	B	142	LEU
1	B	154	ILE
1	C	2	ASN
1	C	19	ILE
1	C	22	PHE

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Mol	Chain	Res	Type
1	C	24	GLN
1	C	30	LEU
1	C	43	GLN
1	C	45	LYS
1	C	50	THR
1	C	59	GLU
1	C	62	LEU
1	C	79	LEU
1	C	118	THR
1	C	136	LYS
1	C	142	LEU
1	C	154	ILE
1	D	2	ASN
1	D	19	ILE
1	D	22	PHE
1	D	24	GLN
1	D	43	GLN
1	D	45	LYS
1	D	50	THR
1	D	59	GLU
1	D	62	LEU
1	D	64	THR
1	D	79	LEU
1	D	118	THR
1	D	136	LYS
1	D	142	LEU
1	D	154	ILE
1	E	2	ASN
1	E	19	ILE
1	E	22	PHE
1	E	24	GLN
1	E	30	LEU
1	E	43	GLN
1	E	45	LYS
1	E	50	THR
1	E	59	GLU
1	E	62	LEU
1	E	79	LEU
1	E	118	THR
1	E	136	LYS
1	E	142	LEU
1	E	154	ILE

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Mol	Chain	Res	Type
1	F	2	ASN
1	F	19	ILE
1	F	24	GLN
1	F	30	LEU
1	F	43	GLN
1	F	45	LYS
1	F	50	THR
1	F	62	LEU
1	F	64	THR
1	F	79	LEU
1	F	118	THR
1	F	136	LYS
1	F	142	LEU
1	F	154	ILE
1	G	2	ASN
1	G	19	ILE
1	G	24	GLN
1	G	30	LEU
1	G	43	GLN
1	G	45	LYS
1	G	50	THR
1	G	59	GLU
1	G	62	LEU
1	G	79	LEU
1	G	118	THR
1	G	136	LYS
1	G	142	LEU
1	G	154	ILE
1	H	2	ASN
1	H	19	ILE
1	H	22	PHE
1	H	24	GLN
1	H	30	LEU
1	H	43	GLN
1	H	45	LYS
1	H	50	THR
1	H	59	GLU
1	H	62	LEU
1	H	64	THR
1	H	79	LEU
1	H	118	THR
1	H	136	LYS

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Mol	Chain	Res	Type
1	H	142	LEU
1	H	154	ILE
1	I	2	ASN
1	I	19	ILE
1	I	24	GLN
1	I	30	LEU
1	I	43	GLN
1	I	45	LYS
1	I	50	THR
1	I	62	LEU
1	I	79	LEU
1	I	118	THR
1	I	136	LYS
1	I	142	LEU
1	I	154	ILE
1	J	2	ASN
1	J	19	ILE
1	J	22	PHE
1	J	24	GLN
1	J	30	LEU
1	J	43	GLN
1	J	45	LYS
1	J	50	THR
1	J	59	GLU
1	J	62	LEU
1	J	64	THR
1	J	79	LEU
1	J	118	THR
1	J	136	LYS
1	J	142	LEU
1	J	154	ILE
1	K	2	ASN
1	K	19	ILE
1	K	24	GLN
1	K	30	LEU
1	K	43	GLN
1	K	45	LYS
1	K	50	THR
1	K	59	GLU
1	K	62	LEU
1	K	64	THR
1	K	79	LEU

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Mol	Chain	Res	Type
1	K	118	THR
1	K	136	LYS
1	K	142	LEU
1	K	154	ILE
1	L	2	ASN
1	L	19	ILE
1	L	22	PHE
1	L	24	GLN
1	L	30	LEU
1	L	43	GLN
1	L	45	LYS
1	L	50	THR
1	L	59	GLU
1	L	62	LEU
1	L	79	LEU
1	L	118	THR
1	L	136	LYS
1	L	142	LEU
1	L	154	ILE
1	M	2	ASN
1	M	19	ILE
1	M	22	PHE
1	M	24	GLN
1	M	30	LEU
1	M	43	GLN
1	M	45	LYS
1	M	50	THR
1	M	59	GLU
1	M	62	LEU
1	M	64	THR
1	M	79	LEU
1	M	118	THR
1	M	136	LYS
1	M	142	LEU
1	M	154	ILE
1	N	2	ASN
1	N	19	ILE
1	N	22	PHE
1	N	24	GLN
1	N	30	LEU
1	N	43	GLN
1	N	45	LYS

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Mol	Chain	Res	Type
1	N	50	THR
1	N	59	GLU
1	N	62	LEU
1	N	79	LEU
1	N	118	THR
1	N	136	LYS
1	N	142	LEU
1	N	154	ILE
1	O	2	ASN
1	O	19	ILE
1	O	22	PHE
1	O	24	GLN
1	O	30	LEU
1	O	43	GLN
1	O	45	LYS
1	O	50	THR
1	O	62	LEU
1	O	79	LEU
1	O	118	THR
1	O	136	LYS
1	O	142	LEU
1	O	154	ILE
1	P	2	ASN
1	P	19	ILE
1	P	22	PHE
1	P	24	GLN
1	P	30	LEU
1	P	43	GLN
1	P	45	LYS
1	P	50	THR
1	P	59	GLU
1	P	62	LEU
1	P	79	LEU
1	P	118	THR
1	P	136	LYS
1	P	142	LEU
1	P	154	ILE
1	Q	2	ASN
1	Q	19	ILE
1	Q	22	PHE
1	Q	24	GLN
1	Q	30	LEU

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Mol	Chain	Res	Type
1	Q	43	GLN
1	Q	45	LYS
1	Q	50	THR
1	Q	59	GLU
1	Q	62	LEU
1	Q	64	THR
1	Q	79	LEU
1	Q	118	THR
1	Q	136	LYS
1	Q	142	LEU
1	Q	154	ILE
1	R	2	ASN
1	R	19	ILE
1	R	24	GLN
1	R	30	LEU
1	R	43	GLN
1	R	45	LYS
1	R	50	THR
1	R	59	GLU
1	R	62	LEU
1	R	79	LEU
1	R	118	THR
1	R	136	LYS
1	R	142	LEU
1	R	154	ILE
1	S	2	ASN
1	S	19	ILE
1	S	24	GLN
1	S	30	LEU
1	S	45	LYS
1	S	50	THR
1	S	59	GLU
1	S	62	LEU
1	S	79	LEU
1	S	118	THR
1	S	136	LYS
1	S	142	LEU
1	S	154	ILE
1	T	2	ASN
1	T	19	ILE
1	T	22	PHE
1	T	24	GLN

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Mol	Chain	Res	Type
1	T	30	LEU
1	T	43	GLN
1	T	45	LYS
1	T	50	THR
1	T	59	GLU
1	T	62	LEU
1	T	79	LEU
1	T	118	THR
1	T	136	LYS
1	T	142	LEU
1	T	154	ILE
1	U	2	ASN
1	U	19	ILE
1	U	24	GLN
1	U	30	LEU
1	U	43	GLN
1	U	45	LYS
1	U	50	THR
1	U	62	LEU
1	U	64	THR
1	U	79	LEU
1	U	116	VAL
1	U	118	THR
1	U	136	LYS
1	U	142	LEU
1	U	154	ILE
1	V	2	ASN
1	V	19	ILE
1	V	22	PHE
1	V	24	GLN
1	V	30	LEU
1	V	43	GLN
1	V	45	LYS
1	V	50	THR
1	V	59	GLU
1	V	62	LEU
1	V	64	THR
1	V	79	LEU
1	V	118	THR
1	V	136	LYS
1	V	142	LEU
1	V	154	ILE

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Mol	Chain	Res	Type
1	W	2	ASN
1	W	19	ILE
1	W	22	PHE
1	W	24	GLN
1	W	30	LEU
1	W	43	GLN
1	W	45	LYS
1	W	50	THR
1	W	62	LEU
1	W	79	LEU
1	W	118	THR
1	W	136	LYS
1	W	142	LEU
1	W	154	ILE
1	X	2	ASN
1	X	19	ILE
1	X	22	PHE
1	X	24	GLN
1	X	30	LEU
1	X	43	GLN
1	X	45	LYS
1	X	50	THR
1	X	59	GLU
1	X	62	LEU
1	X	64	THR
1	X	79	LEU
1	X	118	THR
1	X	136	LYS
1	X	142	LEU
1	X	154	ILE
1	Y	2	ASN
1	Y	19	ILE
1	Y	22	PHE
1	Y	24	GLN
1	Y	30	LEU
1	Y	43	GLN
1	Y	45	LYS
1	Y	50	THR
1	Y	62	LEU
1	Y	79	LEU
1	Y	118	THR
1	Y	136	LYS

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Mol	Chain	Res	Type
1	Y	142	LEU
1	Y	154	ILE
1	Z	2	ASN
1	Z	19	ILE
1	Z	24	GLN
1	Z	30	LEU
1	Z	43	GLN
1	Z	45	LYS
1	Z	50	THR
1	Z	59	GLU
1	Z	62	LEU
1	Z	64	THR
1	Z	79	LEU
1	Z	118	THR
1	Z	136	LYS
1	Z	142	LEU
1	Z	154	ILE
1	1	2	ASN
1	1	19	ILE
1	1	22	PHE
1	1	24	GLN
1	1	30	LEU
1	1	43	GLN
1	1	45	LYS
1	1	50	THR
1	1	59	GLU
1	1	62	LEU
1	1	79	LEU
1	1	118	THR
1	1	136	LYS
1	1	142	LEU
1	1	154	ILE
1	2	2	ASN
1	2	19	ILE
1	2	22	PHE
1	2	24	GLN
1	2	30	LEU
1	2	43	GLN
1	2	45	LYS
1	2	50	THR
1	2	59	GLU
1	2	62	LEU

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Mol	Chain	Res	Type
1	2	79	LEU
1	2	118	THR
1	2	136	LYS
1	2	142	LEU
1	2	154	ILE
1	3	2	ASN
1	3	19	ILE
1	3	22	PHE
1	3	24	GLN
1	3	30	LEU
1	3	43	GLN
1	3	45	LYS
1	3	50	THR
1	3	59	GLU
1	3	62	LEU
1	3	79	LEU
1	3	118	THR
1	3	136	LYS
1	3	142	LEU
1	3	154	ILE
1	4	2	ASN
1	4	19	ILE
1	4	22	PHE
1	4	24	GLN
1	4	30	LEU
1	4	43	GLN
1	4	45	LYS
1	4	50	THR
1	4	59	GLU
1	4	62	LEU
1	4	64	THR
1	4	79	LEU
1	4	118	THR
1	4	136	LYS
1	4	142	LEU
1	4	154	ILE
1	5	2	ASN
1	5	19	ILE
1	5	22	PHE
1	5	24	GLN
1	5	30	LEU
1	5	43	GLN

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Mol	Chain	Res	Type
1	5	45	LYS
1	5	50	THR
1	5	59	GLU
1	5	62	LEU
1	5	79	LEU
1	5	118	THR
1	5	136	LYS
1	5	142	LEU
1	5	154	ILE
1	6	2	ASN
1	6	19	ILE
1	6	22	PHE
1	6	24	GLN
1	6	30	LEU
1	6	43	GLN
1	6	45	LYS
1	6	50	THR
1	6	62	LEU
1	6	79	LEU
1	6	118	THR
1	6	136	LYS
1	6	142	LEU
1	6	154	ILE
1	7	2	ASN
1	7	19	ILE
1	7	24	GLN
1	7	30	LEU
1	7	43	GLN
1	7	45	LYS
1	7	50	THR
1	7	59	GLU
1	7	62	LEU
1	7	79	LEU
1	7	118	THR
1	7	136	LYS
1	7	142	LEU
1	7	154	ILE
1	8	2	ASN
1	8	19	ILE
1	8	24	GLN
1	8	30	LEU
1	8	43	GLN

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Mol	Chain	Res	Type
1	8	45	LYS
1	8	50	THR
1	8	59	GLU
1	8	62	LEU
1	8	64	THR
1	8	79	LEU
1	8	118	THR
1	8	136	LYS
1	8	142	LEU
1	8	154	ILE
1	9	2	ASN
1	9	19	ILE
1	9	22	PHE
1	9	24	GLN
1	9	30	LEU
1	9	43	GLN
1	9	45	LYS
1	9	50	THR
1	9	59	GLU
1	9	62	LEU
1	9	64	THR
1	9	79	LEU
1	9	118	THR
1	9	136	LYS
1	9	142	LEU
1	9	154	ILE
1	a	2	ASN
1	a	19	ILE
1	a	24	GLN
1	a	30	LEU
1	a	43	GLN
1	a	45	LYS
1	a	50	THR
1	a	59	GLU
1	a	62	LEU
1	a	79	LEU
1	a	118	THR
1	a	136	LYS
1	a	142	LEU
1	a	154	ILE
1	b	2	ASN
1	b	19	ILE

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Mol	Chain	Res	Type
1	b	24	GLN
1	b	43	GLN
1	b	45	LYS
1	b	50	THR
1	b	59	GLU
1	b	62	LEU
1	b	79	LEU
1	b	118	THR
1	b	136	LYS
1	b	142	LEU
1	b	154	ILE
1	c	2	ASN
1	c	19	ILE
1	c	22	PHE
1	c	24	GLN
1	c	30	LEU
1	c	43	GLN
1	c	45	LYS
1	c	50	THR
1	c	59	GLU
1	c	62	LEU
1	c	79	LEU
1	c	118	THR
1	c	136	LYS
1	c	142	LEU
1	c	154	ILE
1	d	2	ASN
1	d	19	ILE
1	d	22	PHE
1	d	24	GLN
1	d	30	LEU
1	d	43	GLN
1	d	45	LYS
1	d	50	THR
1	d	59	GLU
1	d	62	LEU
1	d	64	THR
1	d	79	LEU
1	d	118	THR
1	d	136	LYS
1	d	142	LEU
1	d	154	ILE

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Mol	Chain	Res	Type
1	e	2	ASN
1	e	19	ILE
1	e	24	GLN
1	e	30	LEU
1	e	43	GLN
1	e	45	LYS
1	e	50	THR
1	e	59	GLU
1	e	62	LEU
1	e	64	THR
1	e	79	LEU
1	e	116	VAL
1	e	118	THR
1	e	136	LYS
1	e	142	LEU
1	e	154	ILE
1	f	2	ASN
1	f	19	ILE
1	f	24	GLN
1	f	30	LEU
1	f	43	GLN
1	f	45	LYS
1	f	50	THR
1	f	59	GLU
1	f	62	LEU
1	f	64	THR
1	f	79	LEU
1	f	118	THR
1	f	136	LYS
1	f	142	LEU
1	f	154	ILE
1	g	2	ASN
1	g	19	ILE
1	g	22	PHE
1	g	24	GLN
1	g	30	LEU
1	g	43	GLN
1	g	45	LYS
1	g	50	THR
1	g	59	GLU
1	g	62	LEU
1	g	72	LYS

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Mol	Chain	Res	Type
1	g	79	LEU
1	g	118	THR
1	g	136	LYS
1	g	142	LEU
1	g	154	ILE
1	h	2	ASN
1	h	19	ILE
1	h	24	GLN
1	h	30	LEU
1	h	43	GLN
1	h	45	LYS
1	h	50	THR
1	h	62	LEU
1	h	79	LEU
1	h	118	THR
1	h	136	LYS
1	h	142	LEU
1	h	154	ILE
1	i	2	ASN
1	i	19	ILE
1	i	24	GLN
1	i	30	LEU
1	i	43	GLN
1	i	45	LYS
1	i	50	THR
1	i	59	GLU
1	i	62	LEU
1	i	64	THR
1	i	79	LEU
1	i	118	THR
1	i	136	LYS
1	i	142	LEU
1	i	154	ILE
1	j	2	ASN
1	j	19	ILE
1	j	22	PHE
1	j	24	GLN
1	j	30	LEU
1	j	43	GLN
1	j	45	LYS
1	j	50	THR
1	j	59	GLU

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Mol	Chain	Res	Type
1	j	62	LEU
1	j	79	LEU
1	j	118	THR
1	j	136	LYS
1	j	142	LEU
1	j	154	ILE
1	k	2	ASN
1	k	19	ILE
1	k	22	PHE
1	k	24	GLN
1	k	30	LEU
1	k	43	GLN
1	k	45	LYS
1	k	50	THR
1	k	59	GLU
1	k	62	LEU
1	k	64	THR
1	k	79	LEU
1	k	118	THR
1	k	136	LYS
1	k	142	LEU
1	k	154	ILE
1	l	2	ASN
1	l	19	ILE
1	l	22	PHE
1	l	24	GLN
1	l	30	LEU
1	l	43	GLN
1	l	45	LYS
1	l	50	THR
1	l	59	GLU
1	l	62	LEU
1	l	79	LEU
1	l	118	THR
1	l	136	LYS
1	l	142	LEU
1	l	154	ILE
1	m	2	ASN
1	m	19	ILE
1	m	24	GLN
1	m	30	LEU
1	m	43	GLN

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Mol	Chain	Res	Type
1	m	45	LYS
1	m	50	THR
1	m	59	GLU
1	m	62	LEU
1	m	64	THR
1	m	79	LEU
1	m	118	THR
1	m	136	LYS
1	m	142	LEU
1	m	154	ILE
1	n	2	ASN
1	n	19	ILE
1	n	22	PHE
1	n	24	GLN
1	n	30	LEU
1	n	43	GLN
1	n	45	LYS
1	n	50	THR
1	n	59	GLU
1	n	62	LEU
1	n	79	LEU
1	n	118	THR
1	n	136	LYS
1	n	142	LEU
1	n	154	ILE
1	o	2	ASN
1	o	19	ILE
1	o	22	PHE
1	o	24	GLN
1	o	30	LEU
1	o	43	GLN
1	o	45	LYS
1	o	50	THR
1	o	62	LEU
1	o	64	THR
1	o	79	LEU
1	o	118	THR
1	o	136	LYS
1	o	142	LEU
1	o	154	ILE
1	p	2	ASN
1	p	19	ILE

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Mol	Chain	Res	Type
1	p	22	PHE
1	p	24	GLN
1	p	30	LEU
1	p	43	GLN
1	p	45	LYS
1	p	50	THR
1	p	62	LEU
1	p	79	LEU
1	p	118	THR
1	p	136	LYS
1	p	142	LEU
1	p	154	ILE
1	q	2	ASN
1	q	19	ILE
1	q	24	GLN
1	q	30	LEU
1	q	43	GLN
1	q	45	LYS
1	q	50	THR
1	q	59	GLU
1	q	62	LEU
1	q	79	LEU
1	q	118	THR
1	q	136	LYS
1	q	142	LEU
1	q	154	ILE
1	r	2	ASN
1	r	19	ILE
1	r	22	PHE
1	r	24	GLN
1	r	30	LEU
1	r	43	GLN
1	r	45	LYS
1	r	50	THR
1	r	62	LEU
1	r	79	LEU
1	r	118	THR
1	r	136	LYS
1	r	142	LEU
1	r	154	ILE
1	s	2	ASN
1	s	19	ILE

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Mol	Chain	Res	Type
1	s	22	PHE
1	s	24	GLN
1	s	30	LEU
1	s	43	GLN
1	s	45	LYS
1	s	50	THR
1	s	59	GLU
1	s	62	LEU
1	s	79	LEU
1	s	118	THR
1	s	136	LYS
1	s	142	LEU
1	s	154	ILE
1	t	2	ASN
1	t	19	ILE
1	t	22	PHE
1	t	24	GLN
1	t	30	LEU
1	t	43	GLN
1	t	45	LYS
1	t	50	THR
1	t	59	GLU
1	t	62	LEU
1	t	79	LEU
1	t	118	THR
1	t	136	LYS
1	t	142	LEU
1	t	154	ILE
1	u	2	ASN
1	u	19	ILE
1	u	22	PHE
1	u	24	GLN
1	u	30	LEU
1	u	43	GLN
1	u	45	LYS
1	u	50	THR
1	u	59	GLU
1	u	62	LEU
1	u	79	LEU
1	u	118	THR
1	u	136	LYS
1	u	142	LEU

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Mol	Chain	Res	Type
1	u	154	ILE
1	v	2	ASN
1	v	19	ILE
1	v	24	GLN
1	v	30	LEU
1	v	43	GLN
1	v	45	LYS
1	v	50	THR
1	v	62	LEU
1	v	64	THR
1	v	79	LEU
1	v	118	THR
1	v	136	LYS
1	v	142	LEU
1	v	154	ILE
1	w	2	ASN
1	w	19	ILE
1	w	22	PHE
1	w	24	GLN
1	w	30	LEU
1	w	43	GLN
1	w	45	LYS
1	w	50	THR
1	w	59	GLU
1	w	62	LEU
1	w	79	LEU
1	w	118	THR
1	w	136	LYS
1	w	142	LEU
1	w	154	ILE
1	x	2	ASN
1	x	19	ILE
1	x	22	PHE
1	x	24	GLN
1	x	43	GLN
1	x	45	LYS
1	x	50	THR
1	x	59	GLU
1	x	62	LEU
1	x	79	LEU
1	x	118	THR
1	x	136	LYS

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Mol	Chain	Res	Type
1	x	142	LEU
1	x	154	ILE
1	y	2	ASN
1	y	19	ILE
1	y	24	GLN
1	y	30	LEU
1	y	43	GLN
1	y	45	LYS
1	y	50	THR
1	y	59	GLU
1	y	62	LEU
1	y	79	LEU
1	y	118	THR
1	y	136	LYS
1	y	142	LEU
1	y	154	ILE

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (197) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	2	ASN
1	A	7	ASN
1	A	24	GLN
1	A	99	ASN
1	B	2	ASN
1	B	24	GLN
1	B	99	ASN
1	C	2	ASN
1	C	7	ASN
1	C	24	GLN
1	D	2	ASN
1	D	7	ASN
1	D	24	GLN
1	D	99	ASN
1	E	2	ASN
1	E	24	GLN
1	F	2	ASN
1	F	24	GLN
1	F	99	ASN
1	G	2	ASN
1	G	24	GLN
1	G	89	HIS

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Mol	Chain	Res	Type
1	H	2	ASN
1	H	24	GLN
1	H	99	ASN
1	I	2	ASN
1	I	24	GLN
1	I	89	HIS
1	I	99	ASN
1	J	2	ASN
1	J	24	GLN
1	J	99	ASN
1	K	2	ASN
1	K	24	GLN
1	K	99	ASN
1	L	2	ASN
1	L	24	GLN
1	L	89	HIS
1	M	2	ASN
1	M	24	GLN
1	M	99	ASN
1	N	2	ASN
1	N	7	ASN
1	N	24	GLN
1	N	99	ASN
1	O	2	ASN
1	O	7	ASN
1	O	24	GLN
1	O	89	HIS
1	O	99	ASN
1	P	2	ASN
1	P	24	GLN
1	P	99	ASN
1	Q	2	ASN
1	Q	7	ASN
1	Q	24	GLN
1	Q	99	ASN
1	R	2	ASN
1	R	24	GLN
1	R	99	ASN
1	S	2	ASN
1	S	24	GLN
1	S	99	ASN
1	T	2	ASN

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Mol	Chain	Res	Type
1	T	24	GLN
1	T	99	ASN
1	U	2	ASN
1	U	24	GLN
1	U	99	ASN
1	V	2	ASN
1	V	7	ASN
1	V	24	GLN
1	V	99	ASN
1	W	2	ASN
1	W	24	GLN
1	W	99	ASN
1	X	2	ASN
1	X	24	GLN
1	X	99	ASN
1	Y	2	ASN
1	Y	7	ASN
1	Y	24	GLN
1	Y	99	ASN
1	Z	2	ASN
1	Z	24	GLN
1	Z	99	ASN
1	1	2	ASN
1	1	24	GLN
1	1	99	ASN
1	2	2	ASN
1	2	7	ASN
1	2	24	GLN
1	3	2	ASN
1	3	7	ASN
1	3	24	GLN
1	3	99	ASN
1	4	2	ASN
1	4	24	GLN
1	4	99	ASN
1	5	2	ASN
1	5	24	GLN
1	5	99	ASN
1	6	2	ASN
1	6	7	ASN
1	6	24	GLN
1	6	99	ASN

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Mol	Chain	Res	Type
1	7	2	ASN
1	7	7	ASN
1	7	24	GLN
1	7	99	ASN
1	8	2	ASN
1	8	24	GLN
1	8	99	ASN
1	9	2	ASN
1	9	7	ASN
1	9	24	GLN
1	9	99	ASN
1	a	2	ASN
1	a	24	GLN
1	a	99	ASN
1	b	2	ASN
1	b	7	ASN
1	b	24	GLN
1	b	99	ASN
1	c	2	ASN
1	c	24	GLN
1	c	99	ASN
1	d	2	ASN
1	d	24	GLN
1	d	99	ASN
1	e	2	ASN
1	e	24	GLN
1	f	2	ASN
1	f	24	GLN
1	f	99	ASN
1	g	2	ASN
1	g	7	ASN
1	g	24	GLN
1	h	2	ASN
1	h	24	GLN
1	h	99	ASN
1	i	2	ASN
1	i	24	GLN
1	i	99	ASN
1	j	2	ASN
1	j	24	GLN
1	j	99	ASN
1	k	2	ASN

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Mol	Chain	Res	Type
1	k	24	GLN
1	k	99	ASN
1	l	2	ASN
1	l	7	ASN
1	l	24	GLN
1	l	89	HIS
1	l	99	ASN
1	m	2	ASN
1	m	24	GLN
1	m	99	ASN
1	n	2	ASN
1	n	24	GLN
1	n	99	ASN
1	o	2	ASN
1	o	24	GLN
1	o	99	ASN
1	p	2	ASN
1	p	24	GLN
1	q	2	ASN
1	q	7	ASN
1	q	24	GLN
1	q	99	ASN
1	r	2	ASN
1	r	24	GLN
1	r	89	HIS
1	r	99	ASN
1	s	2	ASN
1	s	24	GLN
1	s	99	ASN
1	t	2	ASN
1	t	7	ASN
1	t	24	GLN
1	t	99	ASN
1	u	2	ASN
1	u	24	GLN
1	u	89	HIS
1	u	99	ASN
1	v	2	ASN
1	v	24	GLN
1	v	99	ASN
1	w	2	ASN
1	w	24	GLN

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Mol	Chain	Res	Type
1	w	99	ASN
1	x	2	ASN
1	x	24	GLN
1	x	99	ASN
1	y	2	ASN
1	y	24	GLN
1	y	99	ASN

### 5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

## 5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

## 5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

## 5.6 Ligand geometry [i](#)

79 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the chemical component dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with  $|Z| > 2$  is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z  > 2$	Counts	RMSZ	$\# Z  > 2$
2	SO4	1	157	-	4,4,4	0.21	0	6,6,6	0.18	0
2	SO4	1	158	-	4,4,4	0.20	0	6,6,6	0.20	0
2	SO4	1	159	-	4,4,4	0.18	0	6,6,6	0.26	0
2	SO4	2	157	-	4,4,4	0.34	0	6,6,6	0.58	0
2	SO4	2	158	-	4,4,4	0.13	0	6,6,6	0.34	0
2	SO4	3	157	-	4,4,4	0.20	0	6,6,6	0.23	0
2	SO4	5	157	-	4,4,4	0.15	0	6,6,6	0.29	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
2	SO4	6	157	-	4,4,4	0.22	0	6,6,6	0.29	0
2	SO4	7	157	-	4,4,4	0.13	0	6,6,6	0.27	0
2	SO4	7	158	-	4,4,4	0.15	0	6,6,6	0.23	0
2	SO4	8	157	-	4,4,4	0.11	0	6,6,6	0.14	0
2	SO4	9	157	-	4,4,4	0.12	0	6,6,6	0.31	0
2	SO4	A	157	-	4,4,4	0.25	0	6,6,6	0.27	0
2	SO4	B	157	-	4,4,4	0.13	0	6,6,6	0.37	0
2	SO4	B	158	-	4,4,4	0.18	0	6,6,6	0.26	0
2	SO4	B	159	-	4,4,4	0.18	0	6,6,6	0.31	0
2	SO4	D	157	-	4,4,4	0.11	0	6,6,6	0.33	0
2	SO4	D	158	-	4,4,4	0.16	0	6,6,6	0.29	0
2	SO4	E	157	-	4,4,4	0.17	0	6,6,6	0.27	0
2	SO4	F	157	-	4,4,4	0.30	0	6,6,6	0.40	0
2	SO4	F	158	-	4,4,4	0.14	0	6,6,6	0.10	0
2	SO4	G	157	-	4,4,4	0.23	0	6,6,6	0.23	0
2	SO4	H	157	-	4,4,4	0.13	0	6,6,6	0.66	0
2	SO4	H	158	-	4,4,4	0.08	0	6,6,6	0.24	0
2	SO4	I	157	-	4,4,4	0.15	0	6,6,6	0.44	0
2	SO4	J	157	-	4,4,4	0.19	0	6,6,6	0.23	0
2	SO4	K	157	-	4,4,4	0.24	0	6,6,6	0.31	0
2	SO4	L	157	-	4,4,4	0.21	0	6,6,6	0.17	0
2	SO4	L	158	-	4,4,4	0.35	0	6,6,6	0.32	0
2	SO4	L	159	-	4,4,4	0.13	0	6,6,6	0.14	0
2	SO4	M	157	-	4,4,4	0.17	0	6,6,6	0.47	0
2	SO4	M	158	-	4,4,4	0.16	0	6,6,6	0.20	0
2	SO4	M	159	-	4,4,4	0.16	0	6,6,6	0.21	0
2	SO4	O	157	-	4,4,4	0.18	0	6,6,6	0.14	0
2	SO4	P	157	-	4,4,4	0.24	0	6,6,6	0.25	0
2	SO4	P	158	-	4,4,4	0.16	0	6,6,6	0.21	0
2	SO4	Q	158	-	4,4,4	0.13	0	6,6,6	0.15	0
2	SO4	R	157	-	4,4,4	0.19	0	6,6,6	0.31	0
2	SO4	R	158	-	4,4,4	0.17	0	6,6,6	0.29	0
2	SO4	T	157	-	4,4,4	0.26	0	6,6,6	0.24	0
2	SO4	U	157	-	4,4,4	0.11	0	6,6,6	0.27	0
2	SO4	V	157	-	4,4,4	0.19	0	6,6,6	0.34	0
2	SO4	W	157	-	4,4,4	0.22	0	6,6,6	0.52	0
2	SO4	X	158	-	4,4,4	0.05	0	6,6,6	0.21	0
2	SO4	Y	157	-	4,4,4	0.36	0	6,6,6	0.52	0
2	SO4	Z	157	-	4,4,4	0.19	0	6,6,6	0.19	0
2	SO4	Z	158	-	4,4,4	0.24	0	6,6,6	0.39	0
2	SO4	a	157	-	4,4,4	0.08	0	6,6,6	0.14	0
2	SO4	b	157	-	4,4,4	0.20	0	6,6,6	0.31	0
2	SO4	b	158	-	4,4,4	0.22	0	6,6,6	0.64	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
2	SO4	b	159	-	4,4,4	0.25	0	6,6,6	0.35	0
2	SO4	d	157	-	4,4,4	0.20	0	6,6,6	0.70	0
2	SO4	d	158	-	4,4,4	0.15	0	6,6,6	0.21	0
2	SO4	e	157	-	4,4,4	0.16	0	6,6,6	0.22	0
2	SO4	f	157	-	4,4,4	0.24	0	6,6,6	0.35	0
2	SO4	g	157	-	4,4,4	0.17	0	6,6,6	0.31	0
2	SO4	h	157	-	4,4,4	0.15	0	6,6,6	0.19	0
2	SO4	h	158	-	4,4,4	0.28	0	6,6,6	0.28	0
2	SO4	h	159	-	4,4,4	0.13	0	6,6,6	0.30	0
2	SO4	i	157	-	4,4,4	0.18	0	6,6,6	0.48	0
2	SO4	k	157	-	4,4,4	0.14	0	6,6,6	0.15	0
2	SO4	k	158	-	4,4,4	0.09	0	6,6,6	0.28	0
2	SO4	l	157	-	4,4,4	0.17	0	6,6,6	0.41	0
2	SO4	m	157	-	4,4,4	0.14	0	6,6,6	0.33	0
2	SO4	n	157	-	4,4,4	0.35	0	6,6,6	0.70	0
2	SO4	o	157	-	4,4,4	0.21	0	6,6,6	0.24	0
2	SO4	o	158	-	4,4,4	0.14	0	6,6,6	0.08	0
2	SO4	p	157	-	4,4,4	0.16	0	6,6,6	0.25	0
2	SO4	p	158	-	4,4,4	0.16	0	6,6,6	0.26	0
2	SO4	p	159	-	4,4,4	0.12	0	6,6,6	0.31	0
2	SO4	q	157	-	4,4,4	0.15	0	6,6,6	0.23	0
2	SO4	q	158	-	4,4,4	0.06	0	6,6,6	0.13	0
2	SO4	s	157	-	4,4,4	0.10	0	6,6,6	0.14	0
2	SO4	s	158	-	4,4,4	0.16	0	6,6,6	0.45	0
2	SO4	u	158	-	4,4,4	0.16	0	6,6,6	0.28	0
2	SO4	v	157	-	4,4,4	0.23	0	6,6,6	0.26	0
2	SO4	x	157	-	4,4,4	0.18	0	6,6,6	0.19	0
2	SO4	x	158	-	4,4,4	0.25	0	6,6,6	0.25	0
2	SO4	y	157	-	4,4,4	0.16	0	6,6,6	0.34	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the chemical component dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	SO4	1	157	-	-	0/0/0/0	0/0/0/0
2	SO4	1	158	-	-	0/0/0/0	0/0/0/0
2	SO4	1	159	-	-	0/0/0/0	0/0/0/0
2	SO4	2	157	-	-	0/0/0/0	0/0/0/0
2	SO4	2	158	-	-	0/0/0/0	0/0/0/0
2	SO4	3	157	-	-	0/0/0/0	0/0/0/0
2	SO4	5	157	-	-	0/0/0/0	0/0/0/0

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	SO4	6	157	-	-	0/0/0/0	0/0/0/0
2	SO4	7	157	-	-	0/0/0/0	0/0/0/0
2	SO4	7	158	-	-	0/0/0/0	0/0/0/0
2	SO4	8	157	-	-	0/0/0/0	0/0/0/0
2	SO4	9	157	-	-	0/0/0/0	0/0/0/0
2	SO4	A	157	-	-	0/0/0/0	0/0/0/0
2	SO4	B	157	-	-	0/0/0/0	0/0/0/0
2	SO4	B	158	-	-	0/0/0/0	0/0/0/0
2	SO4	B	159	-	-	0/0/0/0	0/0/0/0
2	SO4	D	157	-	-	0/0/0/0	0/0/0/0
2	SO4	D	158	-	-	0/0/0/0	0/0/0/0
2	SO4	E	157	-	-	0/0/0/0	0/0/0/0
2	SO4	F	157	-	-	0/0/0/0	0/0/0/0
2	SO4	F	158	-	-	0/0/0/0	0/0/0/0
2	SO4	G	157	-	-	0/0/0/0	0/0/0/0
2	SO4	H	157	-	-	0/0/0/0	0/0/0/0
2	SO4	H	158	-	-	0/0/0/0	0/0/0/0
2	SO4	I	157	-	-	0/0/0/0	0/0/0/0
2	SO4	J	157	-	-	0/0/0/0	0/0/0/0
2	SO4	K	157	-	-	0/0/0/0	0/0/0/0
2	SO4	L	157	-	-	0/0/0/0	0/0/0/0
2	SO4	L	158	-	-	0/0/0/0	0/0/0/0
2	SO4	L	159	-	-	0/0/0/0	0/0/0/0
2	SO4	M	157	-	-	0/0/0/0	0/0/0/0
2	SO4	M	158	-	-	0/0/0/0	0/0/0/0
2	SO4	M	159	-	-	0/0/0/0	0/0/0/0
2	SO4	O	157	-	-	0/0/0/0	0/0/0/0
2	SO4	P	157	-	-	0/0/0/0	0/0/0/0
2	SO4	P	158	-	-	0/0/0/0	0/0/0/0
2	SO4	Q	158	-	-	0/0/0/0	0/0/0/0
2	SO4	R	157	-	-	0/0/0/0	0/0/0/0
2	SO4	R	158	-	-	0/0/0/0	0/0/0/0
2	SO4	T	157	-	-	0/0/0/0	0/0/0/0
2	SO4	U	157	-	-	0/0/0/0	0/0/0/0
2	SO4	V	157	-	-	0/0/0/0	0/0/0/0
2	SO4	W	157	-	-	0/0/0/0	0/0/0/0
2	SO4	X	158	-	-	0/0/0/0	0/0/0/0
2	SO4	Y	157	-	-	0/0/0/0	0/0/0/0
2	SO4	Z	157	-	-	0/0/0/0	0/0/0/0
2	SO4	Z	158	-	-	0/0/0/0	0/0/0/0
2	SO4	a	157	-	-	0/0/0/0	0/0/0/0
2	SO4	b	157	-	-	0/0/0/0	0/0/0/0

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	SO4	b	158	-	-	0/0/0/0	0/0/0/0
2	SO4	b	159	-	-	0/0/0/0	0/0/0/0
2	SO4	d	157	-	-	0/0/0/0	0/0/0/0
2	SO4	d	158	-	-	0/0/0/0	0/0/0/0
2	SO4	e	157	-	-	0/0/0/0	0/0/0/0
2	SO4	f	157	-	-	0/0/0/0	0/0/0/0
2	SO4	g	157	-	-	0/0/0/0	0/0/0/0
2	SO4	h	157	-	-	0/0/0/0	0/0/0/0
2	SO4	h	158	-	-	0/0/0/0	0/0/0/0
2	SO4	h	159	-	-	0/0/0/0	0/0/0/0
2	SO4	i	157	-	-	0/0/0/0	0/0/0/0
2	SO4	k	157	-	-	0/0/0/0	0/0/0/0
2	SO4	k	158	-	-	0/0/0/0	0/0/0/0
2	SO4	l	157	-	-	0/0/0/0	0/0/0/0
2	SO4	m	157	-	-	0/0/0/0	0/0/0/0
2	SO4	n	157	-	-	0/0/0/0	0/0/0/0
2	SO4	o	157	-	-	0/0/0/0	0/0/0/0
2	SO4	o	158	-	-	0/0/0/0	0/0/0/0
2	SO4	p	157	-	-	0/0/0/0	0/0/0/0
2	SO4	p	158	-	-	0/0/0/0	0/0/0/0
2	SO4	p	159	-	-	0/0/0/0	0/0/0/0
2	SO4	q	157	-	-	0/0/0/0	0/0/0/0
2	SO4	q	158	-	-	0/0/0/0	0/0/0/0
2	SO4	s	157	-	-	0/0/0/0	0/0/0/0
2	SO4	s	158	-	-	0/0/0/0	0/0/0/0
2	SO4	u	158	-	-	0/0/0/0	0/0/0/0
2	SO4	v	157	-	-	0/0/0/0	0/0/0/0
2	SO4	x	157	-	-	0/0/0/0	0/0/0/0
2	SO4	x	158	-	-	0/0/0/0	0/0/0/0
2	SO4	y	157	-	-	0/0/0/0	0/0/0/0

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

24 monomers are involved in 32 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	3	157	SO4	1	0

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Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	5	157	SO4	2	0
2	6	157	SO4	2	0
2	7	158	SO4	2	0
2	8	157	SO4	1	0
2	A	157	SO4	1	0
2	B	157	SO4	1	0
2	B	158	SO4	3	0
2	B	159	SO4	2	0
2	D	157	SO4	1	0
2	E	157	SO4	1	0
2	F	157	SO4	1	0
2	G	157	SO4	1	0
2	H	157	SO4	1	0
2	H	158	SO4	2	0
2	I	157	SO4	1	0
2	K	157	SO4	1	0
2	L	157	SO4	1	0
2	L	158	SO4	1	0
2	O	157	SO4	1	0
2	P	157	SO4	2	0
2	Q	158	SO4	1	0
2	R	157	SO4	1	0
2	Z	158	SO4	1	0

## 5.7 Other polymers [i](#)

There are no such residues in this entry.

## 5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

## 6 Fit of model and data ⓘ

### 6.1 Protein, DNA and RNA chains ⓘ

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95<sup>th</sup> percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å <sup>2</sup> )	Q<0.9
1	1	154/156 (98%)	-0.52	0 100 100	60, 75, 102, 109	0
1	2	154/156 (98%)	-0.52	0 100 100	59, 76, 103, 111	0
1	3	154/156 (98%)	-0.49	0 100 100	61, 77, 101, 110	0
1	4	154/156 (98%)	-0.46	0 100 100	60, 77, 102, 110	0
1	5	154/156 (98%)	-0.48	0 100 100	62, 75, 102, 112	0
1	6	154/156 (98%)	-0.53	0 100 100	61, 75, 101, 108	0
1	7	154/156 (98%)	-0.55	0 100 100	61, 76, 102, 110	0
1	8	154/156 (98%)	-0.47	0 100 100	63, 77, 100, 112	0
1	9	154/156 (98%)	-0.58	0 100 100	62, 76, 101, 109	0
1	A	154/156 (98%)	-0.51	0 100 100	61, 76, 101, 107	0
1	B	154/156 (98%)	-0.51	0 100 100	61, 77, 102, 107	0
1	C	154/156 (98%)	-0.46	0 100 100	62, 77, 104, 108	0
1	D	154/156 (98%)	-0.56	0 100 100	60, 76, 101, 112	0
1	E	154/156 (98%)	-0.53	0 100 100	61, 76, 100, 108	0
1	F	154/156 (98%)	-0.48	0 100 100	63, 78, 103, 111	0
1	G	154/156 (98%)	-0.48	0 100 100	65, 79, 103, 111	0
1	H	154/156 (98%)	-0.41	0 100 100	63, 79, 103, 110	0
1	I	154/156 (98%)	-0.50	0 100 100	63, 77, 103, 112	0
1	J	154/156 (98%)	-0.51	0 100 100	61, 78, 102, 112	0
1	K	154/156 (98%)	-0.50	0 100 100	59, 77, 102, 110	0
1	L	154/156 (98%)	-0.53	1 (0%) 90 85	64, 78, 102, 110	0
1	M	154/156 (98%)	-0.52	0 100 100	63, 78, 102, 112	0
1	N	154/156 (98%)	-0.52	0 100 100	59, 75, 100, 112	0
1	O	154/156 (98%)	-0.52	0 100 100	58, 75, 101, 107	0

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Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å <sup>2</sup> )	Q<0.9
1	P	154/156 (98%)	-0.44	0 100 100	63, 79, 102, 112	0
1	Q	154/156 (98%)	-0.34	2 (1%) 79 71	64, 79, 105, 114	0
1	R	154/156 (98%)	-0.40	1 (0%) 90 85	62, 78, 104, 110	0
1	S	154/156 (98%)	-0.43	0 100 100	62, 78, 103, 111	0
1	T	154/156 (98%)	-0.46	1 (0%) 90 85	63, 78, 103, 112	0
1	U	154/156 (98%)	-0.42	1 (0%) 90 85	61, 78, 101, 112	0
1	V	154/156 (98%)	-0.44	0 100 100	60, 78, 102, 114	0
1	W	154/156 (98%)	-0.45	0 100 100	60, 78, 102, 113	0
1	X	154/156 (98%)	-0.44	0 100 100	64, 77, 103, 112	0
1	Y	154/156 (98%)	-0.44	3 (1%) 70 61	63, 78, 104, 113	0
1	Z	154/156 (98%)	-0.46	0 100 100	58, 77, 103, 111	0
1	a	154/156 (98%)	-0.42	1 (0%) 90 85	61, 77, 102, 112	0
1	b	154/156 (98%)	-0.52	0 100 100	61, 77, 101, 111	0
1	c	154/156 (98%)	-0.45	0 100 100	62, 78, 102, 109	0
1	d	154/156 (98%)	-0.54	1 (0%) 90 85	62, 79, 104, 112	0
1	e	154/156 (98%)	-0.38	1 (0%) 90 85	60, 78, 103, 112	0
1	f	154/156 (98%)	-0.44	0 100 100	64, 80, 105, 110	0
1	g	154/156 (98%)	-0.47	1 (0%) 90 85	60, 79, 104, 112	0
1	h	154/156 (98%)	-0.45	0 100 100	61, 79, 103, 111	0
1	i	154/156 (98%)	-0.51	0 100 100	64, 77, 103, 110	0
1	j	154/156 (98%)	-0.47	0 100 100	65, 79, 105, 111	0
1	k	154/156 (98%)	-0.43	0 100 100	61, 77, 102, 112	0
1	l	154/156 (98%)	-0.51	0 100 100	60, 77, 101, 111	0
1	m	154/156 (98%)	-0.41	1 (0%) 90 85	62, 78, 103, 109	0
1	n	154/156 (98%)	-0.48	0 100 100	65, 79, 101, 111	0
1	o	154/156 (98%)	-0.37	0 100 100	62, 78, 101, 110	0
1	p	154/156 (98%)	-0.49	0 100 100	61, 77, 102, 112	0
1	q	154/156 (98%)	-0.56	0 100 100	60, 76, 100, 106	0
1	r	154/156 (98%)	-0.40	1 (0%) 90 85	63, 78, 103, 111	0
1	s	154/156 (98%)	-0.40	0 100 100	60, 79, 103, 111	0
1	t	154/156 (98%)	-0.46	0 100 100	59, 79, 103, 112	0

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Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å <sup>2</sup> )	Q<0.9
1	u	154/156 (98%)	-0.47	0 100 100	63, 78, 101, 111	0
1	v	154/156 (98%)	-0.49	0 100 100	63, 77, 102, 110	0
1	w	154/156 (98%)	-0.45	0 100 100	60, 78, 102, 109	0
1	x	154/156 (98%)	-0.41	1 (0%) 90 85	62, 78, 102, 111	0
1	y	154/156 (98%)	-0.43	0 100 100	62, 78, 103, 111	0
All	All	9240/9360 (98%)	-0.47	16 (0%) 95 93	58, 78, 103, 114	0

All (16) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	L	47	ASP	2.8
1	Y	85	GLY	2.7
1	a	48	ASN	2.6
1	g	1	MET	2.6
1	R	48	ASN	2.5
1	U	2	ASN	2.4
1	Y	87	THR	2.4
1	T	107	ASP	2.3
1	e	2	ASN	2.3
1	m	12	ASP	2.2
1	Y	86	GLY	2.2
1	Q	48	ASN	2.1
1	x	48	ASN	2.1
1	d	47	ASP	2.1
1	r	47	ASP	2.0
1	Q	86	GLY	2.0

## 6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

There are no non-standard protein/DNA/RNA residues in this entry.

## 6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

## 6.4 Ligands ⓘ

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(Å <sup>2</sup> )	Q<0.9
2	SO4	L	158	5/5	0.79	0.35	2.83	62,84,145,170	0
2	SO4	u	158	5/5	0.85	0.28	1.24	86,125,128,160	0
2	SO4	T	157	5/5	0.92	0.25	0.10	76,78,121,133	0
2	SO4	h	157	5/5	0.77	0.29	-0.44	100,130,157,175	0
2	SO4	M	157	5/5	0.91	0.17	-0.60	81,92,115,123	0
2	SO4	Z	158	5/5	0.77	0.22	-0.71	85,90,121,137	0
2	SO4	g	157	5/5	0.82	0.20	-0.76	86,106,123,136	0
2	SO4	R	158	5/5	0.93	0.22	-0.96	91,97,120,131	0
2	SO4	E	157	5/5	0.89	0.19	-0.96	71,87,109,118	0
2	SO4	l	157	5/5	0.93	0.20	-1.01	78,90,111,116	0
2	SO4	D	157	5/5	0.94	0.18	-1.02	67,73,117,126	0
2	SO4	K	157	5/5	0.90	0.18	-1.13	42,81,110,119	0
2	SO4	X	158	5/5	0.94	0.17	-1.14	76,97,112,121	0
2	SO4	F	157	5/5	0.93	0.18	-1.20	54,97,117,132	0
2	SO4	v	157	5/5	0.91	0.19	-1.23	77,81,104,104	0
2	SO4	L	157	5/5	0.94	0.13	-1.33	73,101,114,120	0
2	SO4	y	157	5/5	0.94	0.16	-1.39	42,94,108,117	0
2	SO4	o	158	5/5	0.88	0.18	-1.42	107,111,132,147	0
2	SO4	G	157	5/5	0.94	0.19	-1.44	83,88,105,118	0
2	SO4	f	157	5/5	0.95	0.12	-1.45	93,100,107,126	0
2	SO4	h	159	5/5	0.92	0.15	-1.45	91,104,124,157	0
2	SO4	3	157	5/5	0.94	0.14	-1.46	72,108,118,137	0
2	SO4	6	157	5/5	0.95	0.12	-1.47	48,75,123,128	0
2	SO4	q	158	5/5	0.92	0.17	-1.49	81,96,99,121	0
2	SO4	Y	157	5/5	0.92	0.17	-1.55	61,65,132,136	0
2	SO4	R	157	5/5	0.90	0.20	-1.56	84,94,126,131	0
2	SO4	9	157	5/5	0.92	0.16	-1.63	93,94,127,130	0
2	SO4	I	157	5/5	0.91	0.17	-1.68	69,98,122,146	0
2	SO4	s	157	5/5	0.93	0.18	-1.69	92,104,115,129	0
2	SO4	l	157	5/5	0.89	0.17	-1.72	86,105,124,149	0
2	SO4	8	157	5/5	0.95	0.16	-1.74	100,109,110,127	0
2	SO4	m	157	5/5	0.91	0.14	-1.75	63,75,101,108	0
2	SO4	b	157	5/5	0.93	0.18	-1.78	86,93,104,104	0
2	SO4	p	158	5/5	0.90	0.20	-1.78	80,92,115,133	0
2	SO4	o	157	5/5	0.92	0.16	-1.79	62,77,111,127	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
2	SO4	x	157	5/5	0.95	0.13	-2.01	76,83,100,101	0
2	SO4	U	157	5/5	0.94	0.14	-2.08	77,90,114,115	0
2	SO4	5	157	5/5	0.94	0.11	-2.09	56,97,114,129	0
2	SO4	p	157	5/5	0.93	0.14	-2.10	70,91,104,114	0
2	SO4	Z	157	5/5	0.93	0.13	-2.11	81,84,123,131	0
2	SO4	B	157	5/5	0.96	0.10	-2.19	72,77,104,111	0
2	SO4	e	157	5/5	0.93	0.23	-2.23	60,93,120,125	0
2	SO4	i	157	5/5	0.92	0.14	-2.24	70,84,101,104	0
2	SO4	a	157	5/5	0.96	0.13	-2.31	76,83,94,115	0
2	SO4	k	157	5/5	0.95	0.13	-2.50	70,89,123,142	0
2	SO4	B	159	5/5	0.93	0.13	-2.57	67,81,112,118	0
2	SO4	M	159	5/5	0.96	0.10	-2.64	73,81,104,104	0
2	SO4	2	157	5/5	0.95	0.14	-2.68	46,68,103,105	0
2	SO4	s	158	5/5	0.93	0.15	-2.73	71,94,114,119	0
2	SO4	W	157	5/5	0.95	0.14	-2.74	72,75,87,101	0
2	SO4	P	158	5/5	0.94	0.12	-2.80	76,106,106,141	0
2	SO4	x	158	5/5	0.92	0.17	-2.88	78,86,102,112	0
2	SO4	n	157	5/5	0.95	0.12	-2.89	66,79,124,127	0
2	SO4	d	157	5/5	0.91	0.15	-2.93	72,83,98,119	0
2	SO4	b	159	5/5	0.95	0.13	-2.99	64,77,105,115	0
2	SO4	H	157	5/5	0.93	0.16	-3.08	69,77,93,103	0
2	SO4	l	159	5/5	0.95	0.11	-3.18	75,76,103,114	0
2	SO4	A	157	5/5	0.95	0.12	-3.25	53,66,92,103	0
2	SO4	O	157	5/5	0.94	0.13	-3.36	70,88,102,117	0
2	SO4	P	157	5/5	0.97	0.12	-3.39	71,84,110,127	0
2	SO4	J	157	5/5	0.95	0.14	-3.40	54,75,107,118	0
2	SO4	V	157	5/5	0.96	0.10	-3.65	65,85,113,118	0
2	SO4	7	157	5/5	0.96	0.12	-4.05	59,66,96,107	0
2	SO4	b	158	5/5	0.91	0.17	-4.22	83,92,104,129	0
2	SO4	q	157	5/5	0.77	0.27	-	85,129,161,175	0
2	SO4	D	158	5/5	0.77	0.37	-	90,95,140,156	0
2	SO4	M	158	5/5	0.84	0.20	-	96,113,141,164	0
2	SO4	l	158	5/5	0.82	0.36	-	101,114,125,138	0
2	SO4	7	158	5/5	0.82	0.33	-	99,107,149,153	0
2	SO4	F	158	5/5	0.86	0.32	-	113,116,164,178	0
2	SO4	h	158	5/5	0.91	0.30	-	61,90,128,158	0
2	SO4	L	159	5/5	0.93	0.16	-	109,135,141,160	0
2	SO4	p	159	5/5	0.85	0.29	-	90,95,126,139	0
2	SO4	Q	158	5/5	0.90	0.37	-	98,114,139,145	0
2	SO4	B	158	5/5	0.93	0.27	-	73,83,145,150	0
2	SO4	k	158	5/5	0.84	0.26	-	106,108,136,144	0
2	SO4	H	158	5/5	0.83	0.27	-	102,108,130,147	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
2	SO4	d	158	5/5	0.70	0.29	-	99,109,179,188	0
2	SO4	2	158	5/5	0.92	0.20	-	106,114,129,141	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.